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MAGDEP - A Computer Program for Finding Depth

To Basement From Total Field Marine Magnetic Profiles

By

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This report is preliminary and has
not been edited or reviewed for
conformity with Geological Survey
standards and nomenclature

INTRODUCTION

The purpose of the Fortran program MAGDEP is to provide a method for finding the depth to two-dimensional magnetic source bodies when given only the total field magnetic intensity profile. The assumptions made in this depth analysis are that the causitive source body is:

1. A two-dimensional structure
2. Uniformly magnetized
3. A finite or infinite polygon in cross section

Unlike other methods for magnetic depth determinations which assume a given body shape, this technique allows the shape of the source body to emerge during the analysis. The final result is a sequence of spot depths which locate the position and depth to the corners of the magnetic source body. The correct outline of the body is controlled by the shape of the magnetic intensity profile which has been reduced to the pole. This reduced polar profile is also determined by program MAGDEP. Two depth analyses are made by the program; one for thick polygonal bodies and the other for thin (thickness less than depth) or narrow bodies. The determination of the correct magnetic source body shape depends upon a careful interpretation of the results from both depth analyses.

Theoretical Aspects

The method used by program MAGDEP is based on the theoretical work of Nabighian (1972). The reader interested in the details of the theory should refer to his paper ; this section will only outline the theoretical aspects which have been applied to program MAGDEP.

Nabighian (1972) shows that all bodies of polygonal cross section can be obtained by the superposition of a finite number of magnetized steps. He gives the magnetic attraction for such a magnetized step as $\Delta M(x)$ (in eq. 1) and then uses it as the starting point for the analysis. Differentiating the attraction of the magnetized step with respect to x yields the horizontal derivative and this represents the magnetic anomaly due to a thin infinite sheet: $T(x) = \frac{\partial}{\partial x} \Delta M(x)$. The vertical derivative of the magnetized step is obtained by differentiating with respect to y : $T_1(x) = \frac{\partial}{\partial y} \Delta M(x)$. A comparison of the two derivatives shows that $T(x)$ is the negative Hilbert transform of $T_1(x)$. When one forms the complex analytic signal function from the two derivatives: $A(x) = T(x) - i T_1(x)$ many interesting properties arise. Those properties which are used by MAGDEP include:

1. The amplitude $|A(x)|$ is a symmetric function with respect to $x=0$.
2. The complex function $A(x)$ has simple poles at each corner of the polygon.
3. The bell-shaped function $a(x) = |A(x)|^2 = T(x)^2 + T_1(x)^2$ has the property that its half-maximum half-width is equal to the depth to the polygon corner.

From this brief introduction to the theory, the procedure for computing the depths to the magnetic source bodies becomes more defined. In the depth analysis for thick source bodies:

1. The total field intensity profile (residual values) is differentiated with respect to x to yield $T(x)$.
2. A fourier transformation, spectrum modification, and inverse fourier transformation is performed to obtain the complex analytic signal function: $A(x) = T(x) - iT_1(x)$
3. The amplitude square of the signal function is computed to give: $a(x) = T(x)^2 + T_1(x)^2$. The deconvolution of the $a(x)$ curve into a finite number of bell-shaped symmetric curves gives the locations and depths to the corners of the magnetic source body.

The depth analysis for thin (thickness less than depth) magnetic source bodies is accomplished by the same procedure with the exception that in step 1, the total field intensity profile is not differentiated.

In the general case, where the relief of the magnetic source body is irregular, each corner of the body will contribute a symmetric bell-shaped component to the $a(x) = T(x)^2 + T_1(x)^2$ curve. To identify each of these symmetric components, it is necessary to deconvolve the $a(x)$ curve using a nonlinear least square optimization routine. The procedure used by MAGDEP was derived by Marquardt (1963) and is similar to the technique described by Johnson (1969). See Appendix I for more complete details. The amplitude square curve can also be written as: $a(x) = \frac{x^2}{h^2 + (x-c)^2}$ where \propto is proportional to the magnetic susceptibility and the dip of the polygon side, h is the depth to the polygon corner, x is the dependent position variable, and c is

the location of the polygon corner along the profile. Each corner of the magnetic source body will contribute a bell-shaped symmetric curve controlled by the three parameters α , h , and c . At the start of the least square optimazation process, initial values must be assigned to each of the parameters; the 'goodness of fit' of the final solution depends to a large extent upon the quality of the initial guesses.

Program MAGDEP uses the following procedure for locating the symmetric bell-shaped curves and determining the initial values for each of the parameters (no. of parameters = 3 x no. of bell-shaped curves):

1. All local maxima of the $a(x)$ curve are checked to see if:
 - a. The $a(x)$ value is larger than an arbitrary cutoff value.
 - b. The second derivative of the $a(x)$ curve is large enough. (controlled by maximum expected source depth ZMAXKM)
2. If step 1 is true, the initial parameter values are found according to:
 - a. $h(\text{depth})$: find the local half-maximum half-width from the $a(x)$ curve
 - b. α : use the $a(x)$ value at the peak of the bell-shaped curve ($=\alpha^2/h^2$)
 - c. c (position): use the x -position of the peak of the bell-shaped curve

After the initial symmetric curves are located, the least square routine is used to obtain a trial solution. The residual values ($a(x) - \text{computed values}$) are then examined by a similar procedure to determine whether more bell-shaped curves are present. The final solution is the 'best fit' of the $a(x)$ curve by the least squares computed curve (using all bell-shaped curves). Depths and locations of the magnetic source body corners are derived from this final least squares approximation.

Using the Program

Program MAGDEP has been designed so that it can be used in one or two sections depending upon the desires of the user. The first section computes the $a(x)$ curve and the magnetic profile reduced to the pole; the second section does the least squares analysis of the $a(x)$ curve (see theory section). Experience working with MAGDEP has shown that section 2 generally is from 5 to 15 times more time consuming (and expensive) than section 1 and under some circumstances, where high precision is not required, satisfactory results can be obtained by omitting section 2 and using only a visual inspection of the $a(x)$ curve. A second point is that since computation time in the least squares routine increases rapidly with the addition of each bell-shaped curve, the user should consider examining the $a(x)$ curve for each new profile prior to using the least squares routine. Knowing the approximate number of bell-shaped curves present will not only save computation time but will also reduce the chances of obtaining a poor solution because too few bell-shaped curves were specified by the user.* With these notes of warning, the following procedure is suggested for the new user of MAGDEP:

1. Read the theory section in this program description of MAGDEP to become familiar with the terminology used here
2. Look over the program output given in the example (Appendix II)
3. Use the program thru section 1 to obtain the plot of the $a(x)$ curve
4. Examine the $a(x)$ curve to get a feel for the number and type of bell-shaped curves present. The shapes and sizes of the curves present will give a rough estimate of the depths to the body

*Example of computation time for 100 data points on an IBM 360/70 system:
Section 1 only: 4 sec. execution
Section 2 only: thick body analysis only, 7 bell-shaped curves,
4 LSQR iterations: 20 sec. execution

5. Refer to the do's and don'ts section for help in deciding whether the profile should be broken into smaller segments for analysis.
6. Resubmit those parts of the profile which require precision depth estimates using the least squares routine.

Input constants required: (must be right-justified)

Card 1: (required)	Cols: 1-80	Heading (20A4)	
Card 2: (required)	Cols: 1-10	SPDKT	: ship speed in knots (F10.1,2I10)
	11-20	ID	: data interpolation flag =1 want data to be interpolated (must use card 4) =2 use input data as is. (data must be evenly spaced)
	21-30	ISECT	: Number of program sections to use (see above) =1 compute a(x) curve and profile reduced to the pole (optional) =2 do this and use the least squares routine
Card 3: (required)	Cols: 1-10	IBOD	: Types of bodies to do analysis for (I10,F10.1,2I10,F10.3)
	11-20	ZMAXKM	: Maximum estimated depth to the magnetic source body in kilometers
	21-30	IPOLE	: Profile reduced to the pole flag =0 do not want the reduced profile =1 do want the reduced profile (must use Card 5)

Cols: 31-40 INPLT : Request for the initial plots containing the $a(x)$ curve, the initial parameter curve, and the residual total field profile. (Recommended as being useful)
 =0 do not want the plot
 =1 do want the plot

41-50 CUTAX : A factor which is to be multiplied by the internal value of the $a(x)$ curve cutoff point. This is used in determining whether local maxima of the $a(x)$ curve are to be considered as bell-shaped curves. (Recommend using CUTAX = 1. on first run)

Card 4: (use only if ID = 1)
 (2F10.2, F10.1, I10) Cols: 1-10 XSHR : Starting point for the data interpolation in hours of elapsed time from the start of the profile (must be ≥ 0 .)

11-20 XEHR : Ending point for the data interpolation in hours of elapsed time from the start of the profile (can be larger than the time of the last data point)

21-30 DXMIN : New time increment at which data is to be sampled (in minutes) (recommend DXMIN = 2.or 5. for ships magnetic profiles)

31-40 NP : Maximum number of points desired to be in the data set used for the depth analysis (must be less than or equal to 200 - the absolute maximum number of data points allowed)

Card 5: (use only if IPOLE = 1)
 (2F10.1) Cols: 1-10 FINC : Inclination of the earth's magnetic field where the magnetic profile was measured (in degrees)

Cols: 11-20 ANG : Angle between magnetic north and the ship's heading (in degrees where $180 > ANG > 0$)

Card 6: (use only if ISECT = 2)
 (4I10,F10.3) Cols: 1-10 NRES : The number of residual checks which are to be made during the execution of section 2. The normal cycle is: 1st sqr.-residual check - 1st sq - residual check - 1st sq - etc., depending upon NRES. (Recommend initial use of NRES=1)

11-20 NPEAKD : Thick body analysis: The number of bell-shaped curves which are to be used by the least square routine during its first series of iterations. If NPEAKD=0, all bell-shaped curves (max. of 10) with peak values greater than(AXCUT * internal cutoff) will be used. NPEAKD must be less than or equal to 15. (Recommend either: 1. use the two step procedure described above to determine NPEAKD or 2. use NPEAKD=0)

21-30 NPEAKT : Thin body analysis: (same description, requirements, and recommendations as for NPEAKD)

31-40 IOU_T : Flag to control the extra output for the values for:
 1. the least square approximation and 2. the residual values ($a(x) - \text{computed}$)
 =0 no output is desired
 =1 output is desired
 (Recommend initial use of IOU_T=1)

Cols: 41-50 CUTRSD : A factor which is to be multiplied by the internal value of the residual curve cutoff point. This is an arbitrary criteria used in determining whether a local maxima in the residual curve is large enough to allow another bell-shaped curve to be added. (Recommend using CUTRSD=1. on the first run)

Card 7: (use only if ISECT=2)
 (8II0) Cols: 1-10 ITIR(1) : The number of iterations of the least squares routine which are to be made during the first call of MARQL. (MARQL is the name used for the least squares routine) (recommend using ITIR(I) < 3 on first run)

11-20 ITIR(2) : The number of iterations of the least squares routine which are to be made after the first residual check and during the second call of MARQL.

(as required up to ITIR(NRES+1)) : The number of iterations of the least squares routine which are to be made after the NRES residual check and during the NRES+1 call of MARQL.

Card 8 to end of data set: Input of time and total field values
(T60,I2,T6,I3,I2,2F2.0,F5.0) .

6-8 ID : Julian day (sequential starting from 1 on January 1)

9-10 IH : Hour (based on 24 hour day)

11-12 AM : Minute

13-14 AS : Second

Cols: 15-19 TOTMAG : Measured total field intensity value. (The program will subtract a constant field (= average field value) from all measured total field values. If residual field intensity values are used, a minor change will have to be made to the main program (those statements directly following the data interpolation section))

60-61 IY : Year

Note: Data cards of a different format can readily be used by changing statement 100 in Subroutine INPUT to fit the new data format.

Some Do's and Don'ts

At the time of writing of this 'suggestion' section, the program MAGDEP has had limited use; consequently, many of the following suggestions may eventually become obsolete or possibly even misleading. It is hoped that helpful additions and deletions will be made to this description as they are discovered. The current state of the art in computers and algorithms to do non-linear least squares requires a great deal of computing time and memory storage space. It is thus advantageous for the user of MAGDEP and other similar programs, to take a few extra precautions to insure a less expensive and higher quality solution on the first attempt. The two step procedure, as described earlier, is a worthy precaution and consequently a highly recommended DO. Some observations concerning the operation of the least squares routine, which might be useful are:

Least Squares

1. The better the initial guesses to the parameter values, the faster the routine will reach a good solution. It should be noted that the two techniques described by Nabighian (1972) for determining trial depth values (Eqs. 19 and 21) were attempted and found to give adequate results. Unfortunately, occasional bad depth estimates (probably the result of noise in the spectral transformations) were encountered which adversely affected the least square solution. The current method for finding trial depths, although basically simple, has provided more reliable trial values for the least square solution.
2. The least squares routine converges most rapidly for the amplitude of the symmetric curve, less rapidly towards the width of the curve, and least rapidly for the position of the curve.

3. The rapid amplitude convergence also means that smaller amplitude curves adjacent to or on the flanks of large amplitude curves, are often wiped out during the first series of iterations of the least square routine. These small peaks seem to have a better chance for survival after the large peaks are 'locked in'. The residual analysis has been designed to restore these small peaks and other less obvious ones so that they will be included in the depth analysis.
4. To improve a solution when there is a region of small amplitude symmetric curves near an area of large amplitude peaks, it is often best to break the profile into two parts and analyze them separately.
5. Since the number of multiplications for I iterations of the least square routine is approximately equal to $13 \times I \times K \times N$, where K is the number of parameters (i.e., 3 x number of bell-shaped curves) and N is the number of data points, it is advantageous to either:
 - a. Use the largest possible sample interval which does not jeopardize the depth resolution or
 - b. use a smaller sample interval for only a short time period. Both possibilities will minimize the total number of data points.

A list of some other do's and don'ts includes:

<u>DO</u>	<u>Data</u>	<u>DON'T</u>
	Input and Interpolation	
1.	Do modify the subroutine INPUT to accommodate a different data input format if necessary.	1. Don't input residual magnetic values without first changing three statements in the main program.
2.	Do interpolate raw data to the largest sample interval which will not affect the depth resolution. (max resolution is probably about 1/2 the sample interval)	2. Don't use unevenly sampled data without using the data interpolation option.
3.	Do use raw data sampled at the closest interval possible. It is far better to interpolate to a larger sample interval from raw data than to use fewer data cards.	3. Don't use more than 200 input raw data cards or request more than 200 interpolated data points.
4.	Do check the sample listing of the raw data and interpolated data to insure against any input reading errors.	

DO

DON'T

Computing $a(x)$ and
initial deconvolution

1. Do request initial plot on the first run for any new data set (INPLT=1).
 2. Do use a two step procedure and for the first step let the program decide the number of bell-shaped curves to use (let NPEAKT and NPEAKD = 0).
 3. Do consider carefully what the maximum source depth might be ± 2 km (i.e., don't guess wildly and set to 100).
 4. Do examine the plot of the $a(x)$ curve carefully for all symmetric curves which might be detected visually.
 5. Do break up a data profile into smaller segments or change the sample interval if the $a(x)$ curve has groups or areas of high and low amplitude peaks.
1. Don't request more than 15 bell-shaped curves (i.e., don't let NPEAKD or NPEAKT = 16 or more).
 2. Don't request a large number (>7) of bell-shaped curves on a first run, unless the least squares routine will not be used.
 3. Don't set the cutoff for the $a(x)$ curve too low (i.e., AXCUT); otherwise noise becomes a problem.
 4. Don't be afraid to request a plot of the profile reduced to the pole - it's cheap.
 5. Don't worry if the initial parameter curve looks like a poor fit to the $a(x)$ curve - least squares does a good job.

Least squares and
residual analysis

1. Do remember that the time required for each least squares analysis will be proportional to the number of multiplications performed (about $13 \times \# \text{ iter} \times \# \text{ data pts} \times \# \text{ parameters}$).
1. Don't use a cutoff value for the residuals which is too low (i.e., RSDCUT), it may cause a false symmetric curve to be added which may affect the least squares analysis. The internal cutoff has worked all right so far (RSDCUT=1).

2. Do use at least one residual analysis.
3. Do use enough iterations (min.of 3) on the first least square analysis to allow the residual values to develop adequately.
4. Do expect that the least square solution may 'seek' or oscillate through the correct value from iteration to iteration.
2. Don't be discouraged if a seemingly poor fit is obtained after only a few iterations. The program will inform you if major problems arise and will give words of encouragement and advice.

Flow Diagrams

The detailed flow diagrams for the entire program and some of the more important subroutines are included for two reasons:

1. To give the interested (or confused) user an opportunity to look into the 'black box' of MAGDEP
2. To assist in trouble shooting should difficulties arise

For the purpose of illustration, the entire program has been broken into the two sections which were described previously. The physical layout of the program is such that this separation into two sections is not readily apparent. Those subroutines which do not have a detailed flow diagram included here are indicated by dashed boxes; those with flow diagrams are indicated by double solid boxes. The name of the subroutine lies in one corner of the outer most box.

Explanation for flow diagrams:



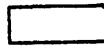
Start and stop



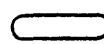
Return from subroutines



Go to



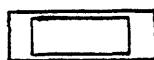
Computational procedure



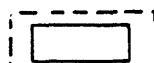
Input and output



Decision

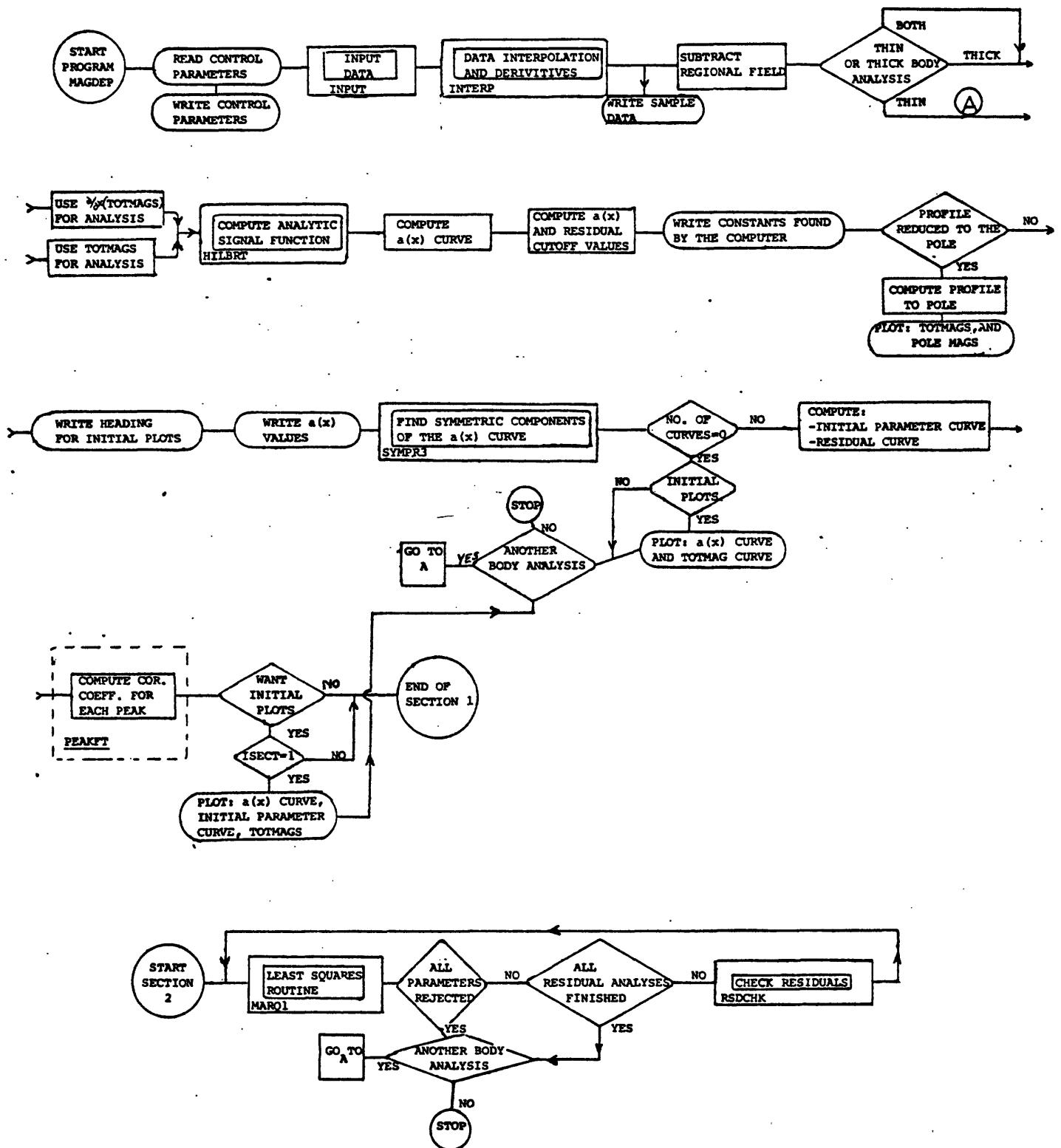


Description of a subroutine which has a detailed flow diagram



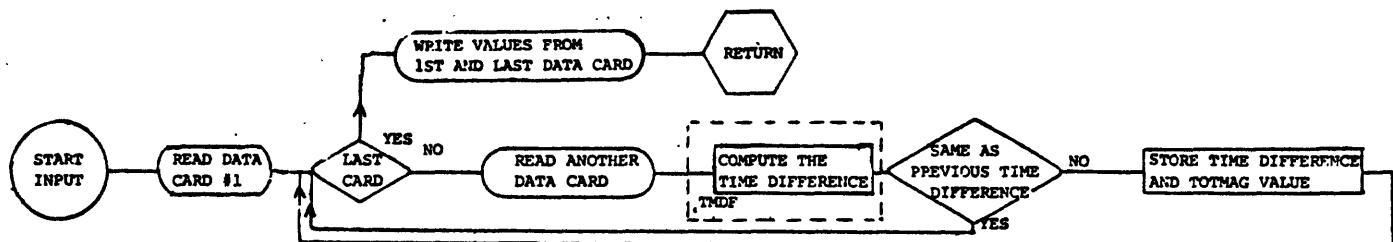
Description of a subroutine which does not have a detailed flow diagram

Flow diagram of entire program MAGDEP:

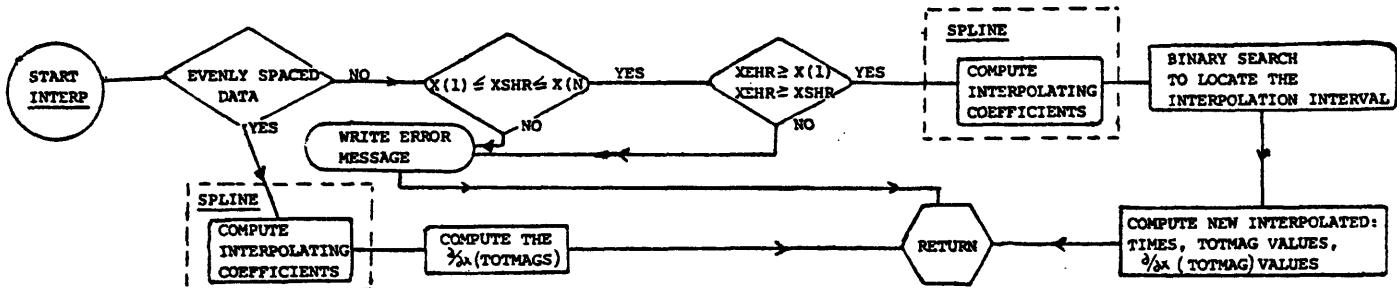


Flow diagrams of important subroutines:

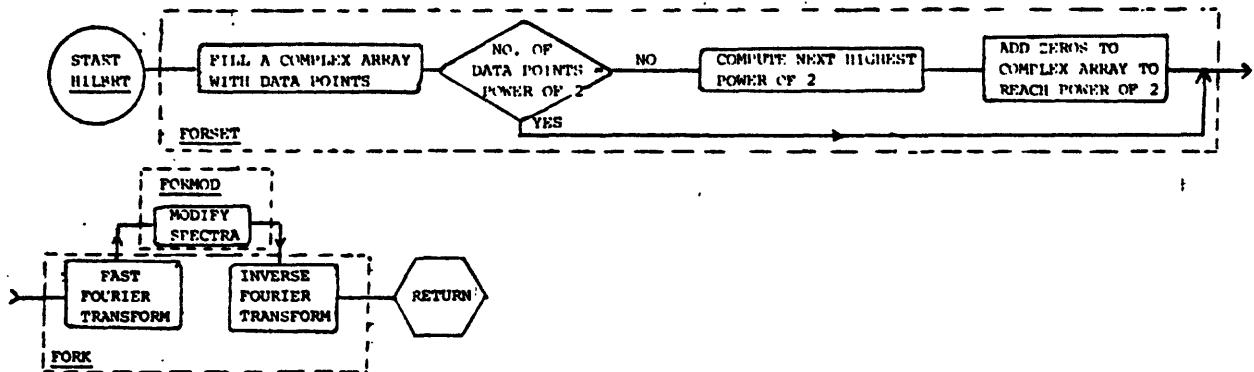
SUBROUTINE INPUT



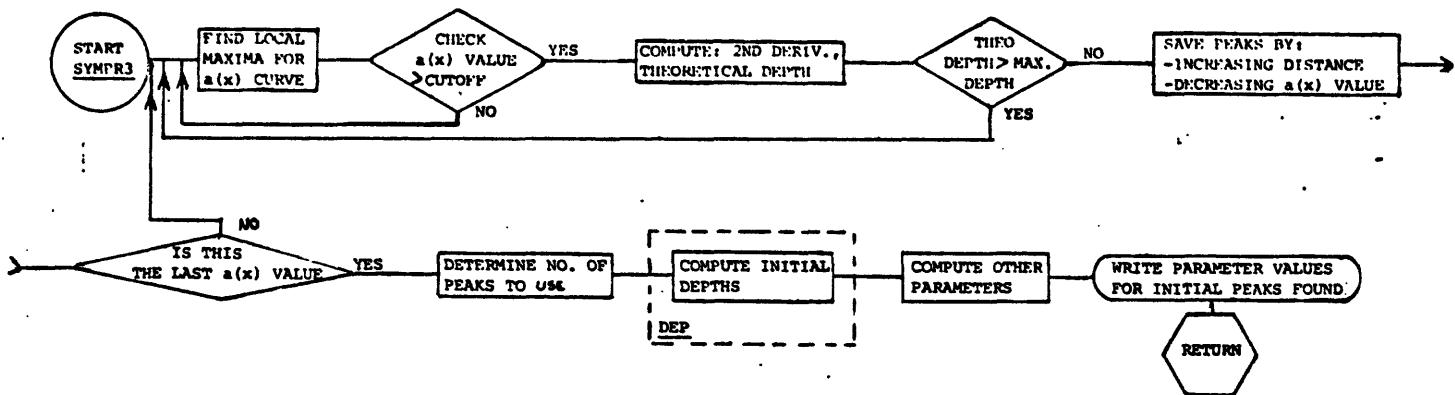
SUBROUTINE INTERP



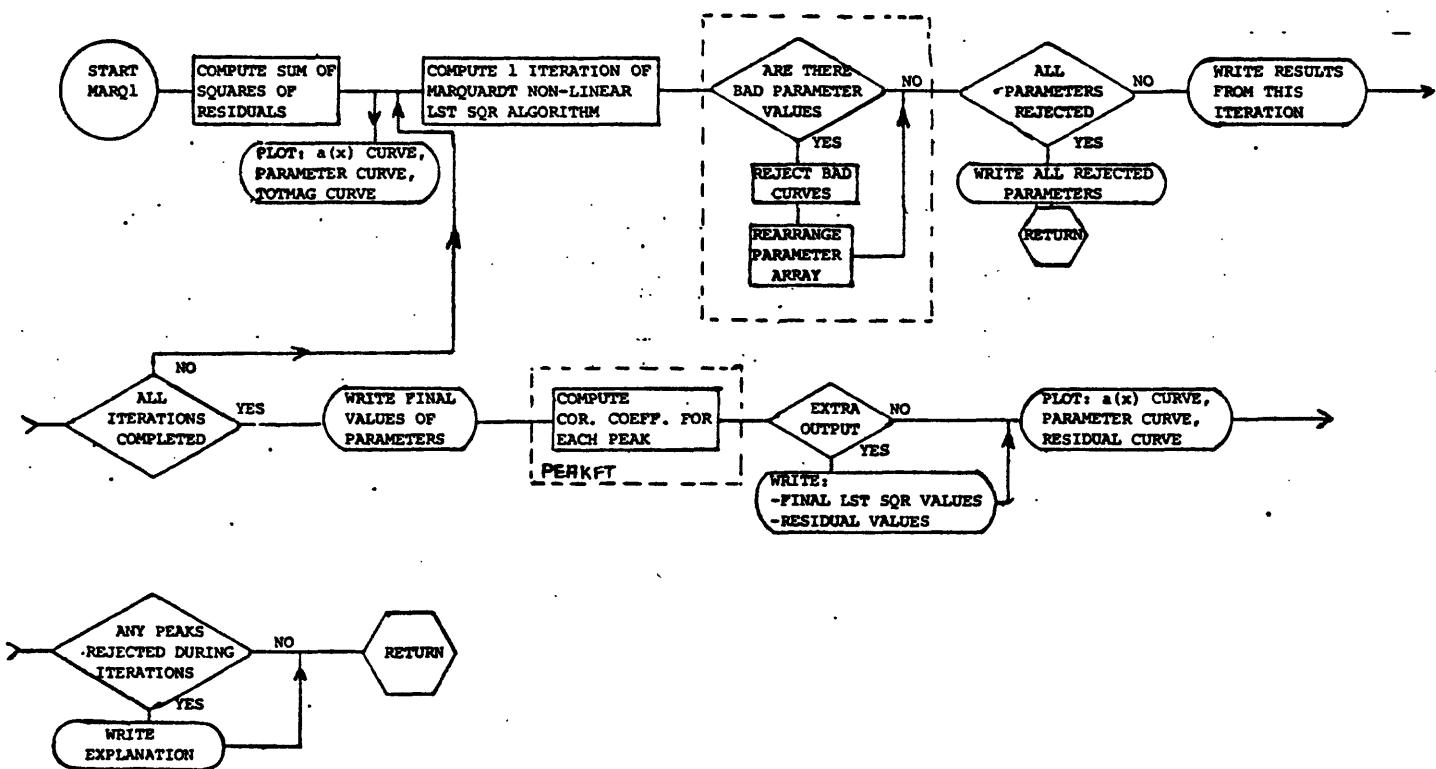
SUBROUTINE HILBRT



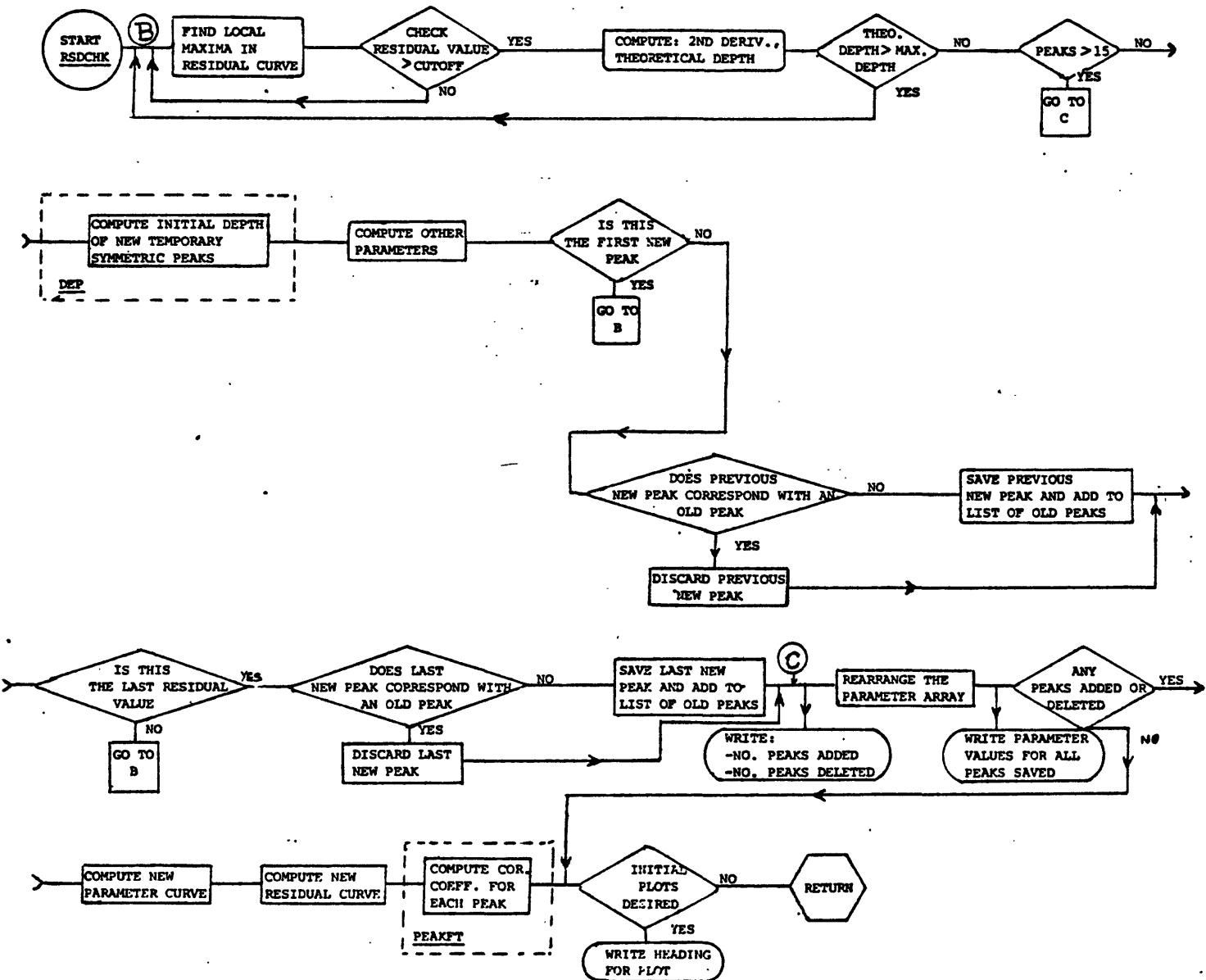
SUBROUTINE SYMPR3



SUBROUTINE MARQ1



SUBROUTINE RSDCHK



Programming Considerations

Listed below are suggestions for making minor changes to program MAGDEP which might be desired at some later time:

A. To change input data format:

1. Modify statement 100 in subroutine INPUT.

B. To use residual magnetic values as input:

1. Modify the statement immediately preceding statement 120 in the main program. Do not subtract AVG from an input (or interpolated) values.

C. To increase the number of data points which may be used, up to N, modify the following statements: Note: wherever N appears, use the desired integer number for N.

1. Main Program:

```
COMMON TIME (N), B(N), C(N), D(N), E(N), F(N)
```

```
COMMON /PRMCRV/XF(N)
```

```
DIMENSION AX(N), TMAG(N), DTMAG(N)
```

```
COMPLEX CAX (next even power of 2 above N)
```

2. Subroutine INTERP

```
COMMON XAXIS(N), B(N), C(N), D(N), E(N), XY(N)
```

3. Subroutine SPLINE

```
REAL*4 H(N), SPP(N)
```

4. Subroutine HILBRT

```
COMPLEX AXM (next even power of 2 above N)
```

5. Subroutine POLRED

```
COMMON X(N), XX(4N)
```

6. Subroutine SYMPR3

COMMON X(N), Q(N), T(N), XX(3N)

7. Function DEP

COMMON X(N), XX(5N)

8. Subroutine PEAKFT

COMMON X(N), Y(N), C(N), D(N), DIFF(N), XX(N)

COMMON /PRMCRV/ F(N)

9. Subroutine PLOT

COMMON X(N), XX(5N)

DIMENSION U(1), V(1), W(1), Y(N,3), SYM(3)

10. Subroutine PRLPLT

DIMENSION SYM(NY), ZY(11), GRAPH(101), Y(N,3), X(1)

11. Subroutine MARQ1

COMMON X(N), TM(N), SS(N), TEM(N), DIFF(N), XX(N)

COMMON /PRMCRV/ F(N)

DIMENSION P(N,45), Y(1)

12. Subroutine DERIV

COMMON X(N), XX(5N)

DIMENSION P(N,45)

13. Subroutine FUNCT

COMMON X(N), XX(5N)

14. Subroutine PRMCHK

COMMON X(N), TT(N), TM(N), XX(3N)

15. Subroutine RSDCHK

COMMON X(N), Q(N), REJ(N), W(N), R(N), F(N)

COMMON /PRMCRV/ FF(N)

16. SUBROUTINE INPUT

IF(N.GT N) Go to 10

D. To increase the number of parameters allowed to KMAX (i.e., 3 times the no. of symmetric curves) change the following: Note: wherever KMAX appears, use the desired integer number for KMAX.

1. Main Program

```
COMMON /SYM/ PARM(KMAX), NPAR
```

2. Subroutine SYMPR3

```
COMMON /SYM/ B(KMAX), K
```

```
DIMENSION AX(1), IDIST (KMAX), IAX(KMAX)
```

```
KMAX='KMAX' (directly below statement 104)
```

3. Subroutine PEAKFT

```
COMMON /SYM/ B(KMAX), K
```

4. Subroutine MARQ1

```
COMMON /SYM/ B(KMAX, K)
```

```
DIMENSION P(200,KMAX)
```

```
DIMENSION GC(KMAX), GS(KMAX), GU(KMAX), D(KMAX), DS(KMAX),
```

```
DG(KMAX), B1(KMAX), A(KMAX, KMAX), AS(KMAX, KMAX)
```

5. Subroutine DERIV

```
COMMON /SYM/ B(KMAX), K
```

```
DIMENSION P(200, KMAX)
```

6. Subroutine PRMCHK

```
COMMON /SYM/ B(KMAX), K
```

8. SUBROUTINE MARQGS

```
DIMENSION A(KMAX, KMHX), B(1), X(1)
```

7. Subroutine RSDCHK

```
COMMON /SYM/ B(KMAX), K
```

```
KMAX='KMAX' (directly after statement 108)
```

E. The following subroutines have been written so that they may be

used outside of MAGDEP; the only changes required are to dimension or common statements:

INTERP (SPLINE)
SPLINE
HILBRT (FORK, FORSET, FORMOD)
FORSET
FORK
FORMOD
CORLAT
MINMAX
PLOT (MINMAX, PRLPLT)
PRLPLT

Note: other subroutines which are called by the main subroutine
are enclosed in parenthesis.

APPENDIX II - Program Listing for MAGDEP

NOTE:

INPUT DATA FORMAT:
NOT BAKTLETT MAIN

AS STORED AT ARMES UNDER TKR15MD 11/28/73

MADE FROM SOURCE DECK COPY 1 OF MAGDEP IN
Coopers Files

DATE = 73332

17/28/02

LEVEL 20

C*****
C -----FORTRAN PROGRAM MAGDEP-----
C FORTRAN IV LEVEL G USGS OPEN FILE 74-1017
C*****
C THE PROGRAM IS DESIGNED TO PROVIDE A METHOD FOR FINDING THE DEPTH TO
C CORNERS OF A TWO-DIMENSIONAL MAGNETIC SOURCE BODY WHEN GIVEN ONLY THE TOTAL
C FIELD MAGNETIC INTENSITY PROFILE. ASSUMPTIONS MADE ABOUT CAUSITIVE SOURCE
C BODY: 1. A TWO-DIMENSIONAL STRUCTURE
C 2. UNIFORMLY MAGNETIZED
C 3. A FINITE OR INFINITE POLYGON IN CROSS SECTION
C TWO DEPTH ANALYSES ARE MADE BY THE PROGRAM: 1. THICK BODIES AND 2. THIN
C BODIES (THICKNESS LESS THAN DEPTH). THE DETERMINATION OF THE CORRECT SOURCE
C BODY SHAPE DEPENDS UPON A CAREFUL INTERPRETATION OF RESULTS FROM BOTH DEPTH
C ANALYSES. THE TECHNIQUES USED BY THE PROGRAM ARE BASED ON THE WORK OF:
C 1. THEORY: NABIGHIAN, MISAC N., GEOPHYSICS, V.37, PP507-517, 1972
C 2. LEAST SQUARES: MARQUARDT, DON W., JOURNAL OF APPLIED MATHEMATICS, V11, P431-441, 1963
C AND JOHNSON, WILLIAM W., GEOPHYSICS, V34, NO1, PP65-74, 1969
C FOR FURTHER DETAILS SEE DOCUMENTATION AVAILABLE FROM OFF. MARINE GEOLOGY
C
C USE OF PROGRAM: THE PROGRAM CAN BE USED IN ONE OR TWO SECTIONS WHICH:
C SECTION 1: COMPUTES A(X) CURVE, MAG PROFILE REDUCED TO POLE, INITIAL
C TRIAL VALUES FOR PARAMETERS OF THE SYMMETRIC CURVES (1 CURVE
C PER BODY CORNER AND 3 PARAMETERS PER CURVE(A,H,C))
C SECTION 2: LEAST SQUARE ANALYSIS, RESIDUAL ANALYSIS
C GENERAL PROCEDURE TO FOLLOW WHEN EXAMINING A NEW DATA PROFILE:
C 1. USE PROGRAM THRU SECT 1 TO SEE PLOT OF A(X) CURVE - USE 5 MIN VALUES
C 2. EXAMINE A(X) CURVE FOR NO. AND TYPES OF SYMMETRIC CURVES. THE HALF-MAX
C HALF-WIDTH IS DEPTH TO BODY CORNER. CHOOSE GOOD VALUE FOR CUTAX.
C 3. BREAK PROFILE INTO SMALLER SEGMENTS: A. TO SEPARATE AREAS OF LARGE AND
C SMALL PEAKS B. TO DELETE OR LOOK CLOSER AT 'FLAT' AREAS C. TO STAY
C WITHIN THE MAX. OF POINTS (INPUT OR INTERPOLATION: MAX=200)
C 4. COMPUTE THRU SECTION 2 USING ITIR(1&2)=SMALL NO. DO ONLY 1 RESIDUAL
C ANALYSIS. AFTER RUN, CHECK VALUE OF CUTRSD
C 5. MAKE FINAL RUN BASED UPON RESULTS FROM STEPS 1 THRU 4
C NOTE: THE EXECUTION TIME FOR THE LEAST SQUARES ROUTINE(SECT 2) IS USUALLY
C FROM 5 TO 15 TIMES GREATER THAN FOR SECT 1. UNDER SOME CIRCUMSTANCES,
C WHERE HIGH PRECISION IS NOT REQUIRED, SATISFACTORY RESULTS CAN BE
C OBTAINED BY OMITTING SECT 2 AND USING ONLY A VISUAL INSPECTION OF
C THE A(X) CURVE.
C INPUT CONSTANTS REQUIRED:
C CARD 1: HEADING FORMAT(20A4) REQUIRED
C CARD 2: REQUIRED FORMAT(F10.1,2I10)
C SPDKT = SHIP SPEED IN KNOTS

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C ID = USER DESIRES DATA INTERPOLATION =1 YES (MUST USE CARD 4)
C =2 NO (DATA MUST BE EVENLY SPACED)
C ISECT = PROGRAM SECTIONS TO USE =1 THRU SECT 1 =2 THRU SECT 2
C CARD 3: IBOD = BODY TYPES TO ANALYSE FOR:
C =1 THICK =2 THIN =3 BOTH
C ZMAXKM = MAX. ESTIMATE DEPTH TO SOURCE BODY IN KM
C IPOLE = WANT PROFILE REDUCED TO POLE: =0 NO =1 YES (USE CARD 5)
C INPLT= WANT INITIAL PLOT (CONTAINS A(X), INITIAL PARAMETER, AND
C RESIDUAL TOTAL FIELD PROFILE): =0 NO =1 YES
C CUTAX = FACTOR WHICH WILL BE MULTIPLIED BY THE INTERNAL VALUE FOR
C THE A(X) CURVE CUTOFF POINT AND USED TO DETERMINE WHETHER
C LOCAL MAXIMA IN A(X) CURVE WILL BE KEPT AS SYMMETRIC CURVE
C (ON FIRST RUN USE CUTAX=1.)
C CARD 4: USE ONLY IF ID=1 FORMAT (3F10.2,110)
C XSHR = START INTERPOLATION--HRS OF ELAPSED TIME FROM START OF LINE
C XEHR = END INTERPOLATION--HRS OF ELAPSED TIME FROM START OF LINE
C DXMIN = TIME INCREMENT AT WHICH DATA IS TO BE INTERPOLATED IN MIN
C NP = MAX. NO. POINTS WANTED IN DATA SET (MUST BE LESS THAN 201)
C CARD 5: USE ONLY IF IPOLE=1 FORMAT (2F10.1)
C FINC = INCLINATION OF EARTHS MAG FIELD WHERE PROFILE WAS MEASURED
C ANG = ANGLE BETWEEN MAG NORTH AND SHIP HEADING (0)ANG) 180 DEG!
C CARD 6: USE ONLY IF ISECT = 2 FORMAT (4F10.3)
C NRES = NO. RESIDUAL CHECKS TO BE MADE (REC NRES=1 ON FIRST RUN)
C NPEAKD = THICK BODY: NO. OF A(X) PEAKS TO BE USED FOR 1ST LST SQRS
C NPEAKT = THIN BODY: NO. OF A(X) PEAKS TO BE USED FOR 1ST LSI SQRS
C ICUT = WANT EXTRA OUTPUT (LST SQR VAL & RESID VAL): =0 NO =1 YES
C CUTRSD = FACTOR USED IN RESIDUAL ANALYSIS (SAME EXPLAN AS CUTAX)
C USE ONLY IF ISECT=2 FORMAT (8I10)
C ITIR(1) = NO. ITERATIONS TO BE MADE DURING 1ST CALL OF LST SQRS
C ITIR(2) = NO. ITERATIONS OF LST SQRS TO BE MADE AFTER RESIDUAL
C NUMBER 1 AND DURING LST SQRS ANALYSIS NUMBER 2
C ITIR(N) = AS REQUIRED UP TO ITIR(NRES+1) = NO. OF ITERATIONS TO BE
C MADE AFTER THE NRES RESIDUAL CHECK AND DURING THE
C NRES+1 CALL OF LST SQRS.
C DATA CARDS: FORMAT (1T60,12,1T6,13,12,2F2.0,F5.0)
C THE REQUIRED INPUT VALUES ON EACH DATA CARD (ACCORDING TO THE FORMAT) ARE:
C IY = YEAR ID = JULIAN DAY IH = HOUR (24 HR DAY) AM = MINUTE
C AS = SECOND TOTMAG = MEASURED TOTAL FIELD INTENSITY VALUE
C NOTE: A MAX. OF 200 DATA CARDS OR 200 INTERPOLATED VALUES ARE ALLOWED
C THE PROGRAM SUBTRACTS A CONSTANT REGIONAL FIELD FROM THE DATA
C DEVELOPED BY: ALAN K COOPER
C OFFICE OF MARINE GEOLOGY

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C U.S. GEOLOGICAL SURVEY
C 345 MIDDLEFIELD ROAD
C MENLO PARK, CALIFORNIA 94025 December 1973
C ****
COMMON TIME(200),B(200),C(200),D(200),E(200),F(200)
COMMON /LSTSQ/XLAM,XU,EPS,OUTPT,SPD,ITER,INPLT
COMMON /RESID/ZMAX,ICALL
COMMON /SYM/PARM(45),NPAR
COMMON /PRMCRV/XF(200)
COMMON /WORK/AA(100),BB(100)
COMMON /CUT/CUTAX,CUTRSD,A MIN,AMAX
COMMON /LSQERR/ERR
DIMENSION AX(200),TMAG(200),ITIR(10)
DIMENSION TITLE(20)
COMPLEX CAX(256)
LOGICAL OUTPT

C -----FORMAT STATEMENTS-----

800 FORMAT(F10.1,2I10)
801 FORMAT(1I0,F10.1,2I10,F10.3)
802 FORMAT(2F10.2,F10.1,I10)
803 FORMAT(2F10.1)
804 FORMAT(4I10,F10.3)
805 FORMAT(8I10)
806 FORMAT(20A4)
900 FORMAT(//,1 VALUES FOR THE A(X) = T(X)**2 + T1(X)**2 CURVE:
\$/ ,10(1,1P5E20.9))
901 FORMAT(1H1) !ESTIMATION OF DEPTH TO THICK MAGNETIC SOURCE BODIES -
\$USING THE HORIZONTAL DERIVITIVE OF THE TOTAL FIELD ,/, ! MAGNETICS
\$ FOR THE ANALYSIS:
902 FORMAT(1H1) !ESTIMATION OF DEPTH TO THIN MAGNETIC SOURCE BODIES - U
\$USING THE TOTAL FIELD MAGNETICS FOR THE ANALYSIS:
903 FORMAT(//,1X,'THE MATRIX IN SUBROUTINE MARQ1 IS SINGULAR!',/
\$! CHECK THE INITIAL VALUES OF THE PARAMETERS OR REFER TO THE MARQ1
\$ PROGRAM DESCRIPTION!)
904 FORMAT(1H1) !PLOT OF A(X) CURVE(+)! INITIAL PARAMETER CURVE(*)! AND
\$ THE RESIDUAL TOTAL FIELD MAGNETIC CURVE(•):!
905 FORMAT(//,1 CORRELATION COEFFICIENT BETWEEN A(X) AND INITIAL PARA
\$ METER CURVE=!,E17.6)
906 FORMAT(1H1) !PLOT OF A(X) CURVE(+) AND THE RESIDUAL TOTAL FIELD MAG
\$NETICS CURVE(*):!
907 FORMAT(//,1 SAMPLE OF INPUT DATA AND INTERPOLATED DATA: (FIRST
\$20 DATA POINTS!),/,! NO. OF INPUT DATA POINTS=!,14,15X,
\$! NO. OF INTERPOLATED DATA POINTS=!,14,/,10X,!TIME!,7X,!TOTALMAGS!,

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```
$27X,'TIME',7X,'TOTMAGS',/(2F15.3,15X,2F15.3)
908 FORMAT(//,' THE DEPTH ANALYSIS FOR THICK BODIES - BASED ON THE HO
$RIZONTAL DERIVITIVE OF THE TOTAL FIELD MAGNETICS - HAS TERMINATED
$SUCCESSFULLY')
909 FORMAT(//,' THE DEPTH ANALYSIS FOR THIN BODIES - BASED ON THE TOT
$AL FIELD MAGNETICS - HAS TERMINATED SUCCESSFULLY')
910 FORMAT(//,' INPUT INFORMATION: /'1' PARAMETERS SUPPLIED BY THE USE
$R:!/,'/,! SHIP SPEED(IKS): SPDKT=!,F6.1,/,! DATA INTERPOLATION: I
$D=!,I4,/,! PROGRAM USE: ISECT=!,I4,/,!
$! BODY TYPES: IBOD=!,I4,/,! MAX.BODY DEPTH(KM): ZMAXKM=!
$F6.1,/,! PROFILE TO POLE: IPOLE='!,I4,I/,! INITIAL PLOTS: INPLT=!
$I4,/,! A(X) CUTOFF: CUTAX=!,F7.3)
911 FORMAT(//,* START INTERPOLATION(HRS): XSHR=!,F7.2,/,!
$! END INTERPOLATION(HRS): XEHR=!,F7.2,/,! SAMPLE INCREMENT(MIN): !
$DXMIN=!,F6.1,/,! MAX.NO.POINTS: NP=!,I6)
912 FORMAT(//,* FIELD INCLINATION(DEG): FINC=!,F7.1,/,! MAG HEADING(DE
$G): ANG=!,F7.1)
913 FORMAT(//,* NO. RESIDUAL CHECKS: NRES=!,I4,/,! NO. PEAKS-THICK BOD
$IES: NPEAKD=!,I4,/,! NO. PEAKS-TINH BODIES: NPEAKT=!,I4,I,/
$! EXTRA OUTPUT: IOUT=!,I4,/,! RESIDUAL CUTOFF: CUTRSD=!,F7.3,
$/,* ITERATIONS: ITIR (NRES+1)=!,I0,I6)
914 FORMAT(1H1,* PLOT OF THE INPUT TOTAL FIELD MAGNETICS(*) AND THE TO
$TAL FIELD MAGNETICS REDUCED TO THE POLE(*): (CONSTANT REGIONAL FI
$ELD REMOVED))
915 FORMAT(1H1,(20A4))
916 FORMAT(//,* CONSTANTS DETERMINED BY THE COMPUTER:!,/,! SHIP SPE
$ED(KM/HR)=!,F6.1,/,! SAMPLE INCREMENT(MIN)=!,F6.1,/,!17X,
$!(HRS)=!,F9.4,/,! CONSTANT REGIONAL FIELD:!,/,! REMOVED FROM INPUT
$DATA=!,F8.0,/,! MAX. BODY DEPTH(HRS)=!,F8.4,/,! A(X) CUTOFF VALUE
$=!,E15.7,/,! RESIDUAL CUTOFF VALUE=!,E15.7)
917 FORMAT(//,* DATA AFTER INTERPOLATION FROM!,F6.2,! TO!,F6.2,! HRS!:!
$)
C-----READ IN PARAMETERS-----
READ(5,806)(TITLE(I),I=1,20)
WRITE(6,915)(TITLE(I),I=1,20)
READ(5,800) SPDKT, ID, ISECT
READ(5,801) IBOD, ZMAXKM, IPOL, INPLT, CUTAX
WRITE(6,910) SPDKT, ID, ISECT, IBOD, ZMAXKM, IPOL, INPLT, CUTAX
IF(ID.EQ.2) GO TO 10
READ(5,802) XSHR, XEHR, DXMIN, NP
WRITE(6,911) XSHR, XEHR, DXMIN, NP
DX=DXMIN/60.
XS=XSHR
```

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XE=XEHR
10 IF(IPOLE.EQ.0) GO TO 20
READ(5,803) FINC,ANG
WRITE(6,912) FINC,ANG
20 IF(ISECT.EQ.1) GO TO 90
READ(5,804) NRES,NPEAKD,NPEAKT,IOUT,CUTRSD
OUTPUT=.TRUE.
IF(IOUT.EQ.0) OUTPT=.FALSE.
KT=NRES+1
READ(5,805) (ITIR(I),I=1,KT)
WRITE(6,913) NRES,NPEAKD,NPEAKT,IOUT,CUTRSD,(ITIR(I),I=1,KT)
90 CONTINUE
C ----- INPUT CONSTANTS -----
IF(ISECT.EQ.1) NPEAKT=0
IF(ISECT.EQ.1) NPEAKD=0
IF(ISECT.EQ.1) CUTRSD=0,
SPD=1.852*SPDKT
ZMAX=ZMAXKM/SPD
IERR=0
IPM=0
EPS=0.15
XU=10.
XLAM=.01
C ----- INPUT DATA -----
CALL INPUT(N,TIME,TMAG)
C ----- INTERPOLATE DATA -----
C ----- ISECT -----
CALL INTERP(N,TIME, TMAG,XS,XE,DX,NI,E,F,DTMAG, ID, NP)
WRITE(6,917) XSHR,XEHR,IFED,EQ4) Go To 70
C ----- INSET -----
CALL PRINT(E(1),E(NI)) TO CONTINUE
WRITE(6,907) N,NI,(TIME(I),TMAG(I),E(I),F(I),I=1,20)
N=NI
CALL MINMAX(N,F,TMIN,TMAX)
AVG=(TMAX+TMIN)/2.
DO 120 I=1,N
TIME(I)=E(I)
TMAG(I)=F(I)-AVG
120 CONTINUE
L=1
IF(IBOD.EQ.3) LE=2
DO 1001 JK=1,LE
ICAL=L
ITK=1
KBOD=IBOD+JK

```

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```
C -----COMPUTE A(X) CURVE-----
      GO TO (200,200,201,200,201), KBOD
  200 CALL HILBRT(N,NC,DTMAG,CAX)
  201 CALL HILBRT(N,NC,TMAG,CAX)
  210 CONTINUE
      DO 211 I=1,N
  211 AX(I)=CABS(CAX(I))*#2
C -----COMPUTE CUTOFF POINT FOR A(X) AND RESIDUAL CURVES-----
      IF(JK.EQ.2) GO TO 212
      CALL MINMAX(N,AX,AMIN,AMAX)
      CUTOFF=SQRT(AMAX*AMIN)
      CUTAX=CUTOFF*CUTAX
      CUTRSRSD=CUTOFF*CUTRSRSD
      IF(IPD.EQ.2) DX=TIMEC(2)-TIMEC(1)
      IF(IPD.EQ.2) DXMIN=DX,DX
      ----OUTPUT COMPUTED CONSTANTS-----
      WRITE(6,916) SPD,DXMIN,DX,Avg,ZMAX,CUTAX,CUTRSRSD
  212 CONTINUE
C -----MAGNETIC PROFILE REDUCED TO THE POLE-----
      IF(IPOLE.EQ.0.OR.IPM.NE.0) GO TO 228
      CR=.01745329
      BINC=ATAN(TAN(IFINC*CR)/ABS(COS(ANG*CR)))/CR
      GO TO 222
  221 BINC=90.
  222 PHI2=180.-2.*BINC
  225 CALL POLRED(N,CAX,PHI2)
      WRITE(6,914)
      CALL PLOT(N,2,TMAG,B,B)
  228 IPM=1
C -----FIND SYMMETRIC COMPONENTS OF A(X)-----
      GO TO (231,231,232,231,232), KBOD
  231 WRITE(6,901)
      WRITE(6,900) (AX(I),I=1,N)
      CALL SYMPR3(N,NPEAKD,AX)
      GO TO 230
  232 WRITE(6,902)
      WRITE(6,900) (AX(I),I=1,N)
      AX(I)=AX(2)
      CALL SYMPR3(N,NPEAKT,AX)
  230 CONTINUE
      IF(NPAR.EQ.0) GO TO 233
      CALL FUNCT(PARM,NPAR,XF,N)
      DO 234 JT=1,N
```

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```
234 E(JT)=AX(JT)-XF(JT)
      CALL PEAKFT(N,AX)
C      -----PLOT AX()! TOTMAGS! & INITIAL PARAMETER CURVE-----
      233 IF(INPLT.EQ.0) GO TO 250
      CALL MINMAX(N,AX,A MIN,A MAX)
      CALL SETUP(N,TMAG,B,A MAX,A MIN)
      IF(NPAR.EQ.0) GO TO 235
      WRITE(6,904)
      IF(I SECT.GE.2) GO TO 252
      CALL FUNCT(PARM,NPAR,C,N)
      CALL CORLAT(N,AX,C,VAL)
      WRITE(6,905) VAL
      GO TO 237
      235 WRITE(6,906)
      CALL PLOT(N,2,AX,B,AX)
      GO TO 1000
      237 CALL PLOT(N,3,AX,C,B)
      GO TO 1000
      250 IF(I SECT.LT.2) GO TO 1000
      252 CONTINUE
C      -----LEAST SQUARES' ROUTINE-----
      ITER=ITIR(ITK)
      300 CALL MARQIN(AX,$310)
      IF(NPAR.EQ.0) GO TO 1000
      GO TO 320
      310 WRITE(6,903)
      GO TO 1000
      320 IF(NRES-ITK.LT.0) GO TO 1000
      ITK=ITK+1
C      -----CHECK RESIDUALS-----
      CALL RSDCHK(N,AX)
      GO TO 252
      1000 CONTINUE
      GO TO (400,400,401,400,401),KBOD
      400 WRITE(6,908)
      GO TO 1001
      401 WRITE(6,909)
      1001 CONTINUE
      STOP
      END
```

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INTERP

```
SUBROUTINE INTERP(N,X,Y,XSTART,XSTOP,XINC,N1,X1,Y1,DY1,IDL,NP)
C THIS SUBROUTINE WILL INTERPOLATE THE VALUES FROM ARRAY Y(N) FROM XSTART
C TO XSTOP AT INCREMENTS OF XINC. THREE ARRAYS ARE RETURNED :
C      X1(N1) = REAL ARRAY CONTAINING THE INTERPOLATED X VALUES
C      Y1(N1) = REAL ARRAY CONTAINING THE EVENLY SPACED INTERPOLATED VALUE
C      DY1(N1)= REAL ARRAY CONTAINING THE DERIVATIVES OF Y(N) AT EVENLY
C           SPACED INCREMENTS XINC (IE AT SAME LOCATIONS AS Y1(N1))

C THE OTHER VARIABLES INCLUDE:
C      X(N) = REAL ARRAY CONTAINING THE INITIAL VALUES ALONG THE X-AXIS
C      Y(N) = REAL ARRAY CONTAINING THE INITIAL VALUES TO BE INTERPOLATED
C      XSTART = FIRST X VALUE WHERE THE INTERPOLATION IS TO START
C      XSTOP = LAST X VALUE FOR THE INTERPOLATION
C      XINC = INCREMENT OF X AT WHICH EVENLY SPACED DATA WILL BE RETURNED
C      N = NUMBER OF DATA POINTS IN THE INPUT ARRAY Y(N)
C      N1 = NUMBER OF DATA POINTS IN THE OUTPUT ARRAY Y1(N1) AND DY1(N1)
C      NP = MAXIMUM NUMBER OF POINTS ALLOWED IN OUTPUT ARRAYS
C      ID = IDENTIFICATION MARKER TO INSTRUCT THE SUBROUTINE, THE VALUES:
C           =0 : DO AN INTERPOLATION USING ALL PARAMETERS, THREE ARRAYS
C           ARE RETURNED, AND SUBROUTINE SPLINE IS NOT CALLED
C           =1 : SAME AS FOR 0 EXCEPT SUBROUTINE SPLINE IS CALLED
C           =2 : THE SUBROUTINE WILL ONLY RETURN THE DERIVATIVE OF THE
C           INPUT CURVE AT EACH OF THE EVENLY SPACED X VALUES
C THIS IS A GENERAL SUBROUTINE FOR USE OUTSIDE OF THIS PROGRAM
C MAXIMUM COMMON USAGE: XAXIS(0),B(0),C(0),D(0),E(0),XY(0)
C COMMON XAXIS(200),B(200),C(200),D(200),E(200),XY(200)
C DIMENSION X(1),Y(1),X1(1),Y1(1),DY1(1)
100 FORMAT(//1X,'XSTART IS OUT OF THE RANGE OF THE INPUT DATA')
101 FORMAT(//1X,'XSTOP IS OUT OF THE RANGE OF THE INPUT DATA')
IEND=0
IF(ID.EQ.2) GO TO 15
C -----CHECK INTERPOLATION RANGE-----
1 IF(XSTART.LE.XIN) AND .XSTART.GE.X(1)) GO TO 8
WRITE(6,100)
IEND=1
8 IF(XSTOP.GE.X(1).AND.XSTOP.GE.X(1)) GO TO 9
WRITE(6,101)
RETURN
9 IF(IEND.EQ.1) RETURN
IF(ID.EQ.0) GO TO 1
C -----COMPUTE THE INTERPOLATING CONSTANTS-----
CALL SPLINE(N,X,Y,B,C,D)
C -----BINARY SEARCH-----
```

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```
1  XX=XSTART
  I=1
  J=N
  2  IF(J-1.EQ.1) GO TO 6
  K=(I+J)/2
  IF(XX-X(K)) .3,.5,.4
  3  J=K
  IF(X(K)-XX,GT,X(K)-X(K-1)) GO TO 2
  J=K-1
  GO TO 7
  4  I=K
  IF(XX-X(K),GT,X(K+1)-X(K)) GO TO 2
  J=K
  GO TO 7
  5  J=K
  GO TO 7
  6  J=1
  7  IF(XSTOP,GT,X(N)) XSTOP=X(N)
  M=(XSTOP-XSTART)/XINC+1
  IF(M.GT.NP) M=NP
  C-----DO INTERPOLATION - COMPUTE DERIVITIVES-----
  11  DO 10 I=1,M
  12  IF(XX.LT.X(J+1)) GO TO 13
  IF(J+1.NE.N) GO TO 14
  N1=I-1
  RETURN
  14  J=J+1
  GO TO 12
  13  IF(XX,GT,XSTOP) RETURN
  N1=I
  DX=XX-X(J)
  Y1(I)=Y(J)+DX*(B(J)+DX*(C(J)+DX*D(J)))
  DY1(I)=B(J)+DX*(2.*C(J)+3.*DX*D(J))
  X1(I)=XX
  XX=XX+XINC
  10  CONTINUE
  RETURN
  C-----ONLY COMPUTE DERIVITIVES-----
  15  CALL SPLINE(N,X,Y,B,C,D)
  KK=N-1
  N1=N
  DO 16 I=1,N
```

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INTERP

X1(I)=X(I)

16 Y1(I)=Y(I)

DO 17 I=1, KK

17 DY1(I)=B(I)

DY1(N)=B(KK)

20 RETURN

END

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SPLINE

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```
      SUBROUTINE SPLINE(N,X,Y,B,C,D)
C THIS SUBROUTINE COMPUTES THE SPLINE COEFFICIENTS NECESSARY TO DO DATA
C INTERPOLATION. THE INPUT PARAMETERS ARE:
C   N = NO. OF DATA POINTS
C   X(N) = REAL ARRAY CONTAINING THE ABSISSA OF THE DATA POINTS
C   Y(N) = REAL ARRAY CONTAINING THE ORDINATES OF THE DATA POINTS
C   B(N-1), C(N-1), D(N-1) = REAL ARRAYS CONTAINING THE SPLINE
C COEFFICIENTS
C
C   REAL*4 X(1), Y(1), B(1), C(1), D(1)
C
C   REAL*8 DELTA,DELTAI
C   REAL*4 H(200),SPP(200)
C
C   NM1 = N-1
C   DO 10 I=1,NM1
C     H(I) = X(I+1) - X(I)
C     DELTA = (Y(I+1) - Y(I))/H(I)
C     IF (I.NE.1) B(I) = DELTA - DELTA
C     DELTA = DELTA
C 10  CONTINUE
C   -----FORWARD ELIMINATION-----
C   C(2) = 2.0D0*(H(1) + H(2))
C   DO 20 I = 3,NM1
C     C(I) = 2.0D0*(X(I+1) - X(I-1)) - H(I-1)**2/C(I-1)
C     B(I) = B(I) - H(I-1)*B(I-1)/C(I-1)
C 20  CONTINUE
C   -----BACK SUBSTITUTION-----
C   SPP(1) = 0.0D0
C   SPP(N) = 0.0D0
C   SPP(N-1) = B(N-1)/C(N-1)
C   NM2 = N-2
C   DO 30 I = 2,NM2
C     J = N-I
C     SPP(J) = (B(J) - H(J)*SPP(J+1))/C(J)
C 30  CONTINUE
C
C   DO 40 I = 1,NM1
C     B(I) = (Y(I+1) - Y(I))/H(I) - (SPP(I+1) + 2.0D0*SPP(I))*H(I)
C     C(I) = 3.0D0*SPP(I)
C     D(I) = (SPP(I+1) - SPP(I))/H(I)
C 40  CONTINUE
C   RETURN
C   END
```

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SUBROUTINE FORSET(N,F,NC,CF)

C THIS SUBROUTINE PREPARES THE DATA SET F(N) FOR ENTRY INTO THE FAST

C FOURIER TRANSFORM SUBROUTINE FOK. NC IS THE NUMBER OF POINTS IN THE

C MODIFIED DATA SET IF ZEROS ARE ADDED TO THE NEAREST POWER OF 2. CF(1)

C IS A COMPLEX ARRAY WHICH CONTAINS THE DATA READY FOR INPUT INTO FOK.

DIMENSION F(1)

COMPLEX CF(1)

C -----FIND NEXT HIGHEST POWER OF 2-----

IR=ALOG(1.*N)/.69315

NC=2***(IR+1)

NX=N+1

C -----FILL THE COMPLEX ARRAY-----

DO 1 I=1,N

CF(I)=F(I)

1 CONTINUE

IF(N.EQ.NC) RETURN

C -----ADD ZEROS-----

DO 2 I=NX,NC

CF(I)=0.

2 CONTINUE

RETURN

END

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HILBRT

SUBROUTINE HILBRT(N,NC,FX,FXM)

C THIS SUBROUTINE WILL COMPUTE THE COMPLEX ARRAY AX(1:NC) = T(X) - IT1(X)
C FROM A REAL ARRAY FX(1:N). THE REAL ARRAY IS FOUR. TRANS. MODIFIED, AND
C INVERSE FOUR. TRANS. SUCH THAT: THE REAL PART OF THE COMPLEX OUTPUT ARRAY
C EQUALS THE INPUT ARRAY, AND THE IMAGINARY PART OF THE OUTPUT ARRAY IS THE
C HILBERT TRANSFORM OF THE INPUT ARRAY
C FX(1:N) = A 1-DIMENSIONAL REAL ARRAY CONTAINING THE DATA TO BE TRANSFORMED
C FXM(1:NC) = 1-DIMENSIONAL COMPLEX ARRAY CONTAINING THE TRANSFORM

DIMENSION FX(1:1)

COMPLEX FXM(1:1),AXM(1:256)

C -----FAST FOURIER TRANSFORM-----

CALL FORSET(N,FX,NC,AXM)

CALL FORK(NC,AXM,-1.0)

C -----MODIFY SPECTRUM-----

CALL FORMOD(NC,AXM,FXM)

C -----INVERSE FAST FOURIER TRANSFORM-----

CALL FORK(NC,FXM,1.0)

RETURN

END

G LEVEL 20 FORK

DATE = 73332

17/28/02

SUBROUTINE FORKLX,CX,SIGNI

C FAST FOURIER TRANSFORM ROUTINE WRITTEN BY JON CLARBOUT, GEOPHYSICS
C DEPARTMENT, STANFORD UNIVERSITY, 1970.

```
C
C   CX(K) = SQRT(1/LX) SUM (CX(J)*EXP(2*PI*SIGNI*(J-1)*(K-1)/LX))  
C   J=1
C
C   COMPLEX CX(LX),CARG,CEXP,CW,CTEMP
C   J=1
C   SC=SQRT(1./LX)
DO 5 I=1,LX
1 IF(I.GT.J)GO TO 2
CTEMP=CX(J)*SC
CX(J)=CX(I)*SC
CX(I)=CTEMP
2 M=LX/2
3 IF(J.LE.M)GO TO 5
J=J-M
M=M/2
IF(M.GE.1)GO TO 3
5 J=J+M
L=1
6 ISTEP=2*L
DO 8 M=1,L
CARG=(0.,1.)*(3.14159265*SIGNI*(M-1))/L
CW=CEXP(CARG)
DO 8 I=M,LX,ISTEP
CTEMP=CW*CX(I+L)
CX(I+L)=CX(I)-CTEMP
8 CX(I)=CX(I)+CTEMP
L=ISTEP
1 IF(L.LT.LX)GO TO 6
9 RETURN
END
```

LEVEL 20

POLRED

DATE = 73332

17/28/02

SUBROUTINE POLRED(N,A,PHI2)

C THIS SUBROUTINE TAKES THE COMPLEX ANALYTIC FUNCTION: A(X)=T(X)-IT(X)
C AND USES IT TO OBTAIN A MAGNETIC PROFILE WHICH HAS BEEN REDUCED TO THE
C POLE. THE PROFILE OUTPUT IS STORED IN ARRAY B

C MAXIMUM COMMON USAGE: X(N),B(N),XX(0)

COMMON X(200),B(200),XX(800)

COMPLEX A(1)

DX=X(2)-X(1)

CON=3.141597/180.

PHI2=PHI2*CON

R1=COS(PHI2)

R2=SIN(PHI2)

POLMAG=0.

WRITE(6,100)

100 FORMAT(10 INPUT DATA REDUCED TO THE POLE.)

DO 1 I=1,N

POLMAG=POLMAG+(REAL(A(I))*R1-AIMAG(A(I))*R2)*DX

B(I)=POLMAG

WRITE(6,200) I,B(I)

200 FORMAT(5X,I3,5X,F12.5)

1 CONTINUE

RETURN

END

/

LEVEL 20 FORMOD

DATE = 73332

17/28/02

SUBROUTINE FORMOD(INC,CF,CFM)

C THIS SUBROUTINE WILL MODIFY THE FOURIER TRANSFORM INTO A HILBERT
C TRANSFORM BY MAKING THE FOLLOWING MODIFICATIONS:

H(W)=F(W) FOR W=0

H(W)=2F(W) FOR F(W)>0

H(W)=0 FOR F(W)<0

NC IS THE NUMBER OF POINTS IN CF(INC)

C THE ORIGINAL ARRAY CF(INC) IS SAVED AND THE MODIFIED TRANSFORM ARRAY CFM(INC)
C IS RETURNED

COMPLEX CF(1),CFM(1)

NB=NC/2+1

NA=NB+1

CFM(1)=CF(1)

DO 1 I=2,NB

CFM(I)=2.*CF(I)

1 CONTINUE

DO 2 I=NA,NC

CFM(I)=0,

2 CONTINUE ~~CFM(NB)=0~~

RETURN
END

SUBROUTINE SYMPR3(N,IPAK,A(X))

C THIS SUBROUTINE FINDS THE LOCATION OF EACH SYMMETRIC COMPONENT OF THE A(X) CURVE BY THE FOLLOWING CRITERIA: 1. FIND THE LOCAL MAXIMA OF A(X) - TEST ADJACENT POINTS. 2. CHECK THE 2ND DERIVATIVE AT EACH MAXIMA - A THEORETICAL DEPTH BASED ON A CONTINUOUS 2ND DERIVATIVE IS COMPUTED AND THIS MUST BE LESS THAN ZMAX. THE SUBROUTINE COMPUTES THE PARAMETERS FOR A NUMBER OF THESE SYMMETRIC COMPONENTS WHICH IS DETERMINED BY:
 C 1. THE USER SPECIFYING A VALUE FOR IPAK, OR
 C 2. SETTING IPAK=0 WHICH USES ALL COMPONENTS WHICH HAVE AN A(X) VALUE LARGER THAN THE A(X) CUTOFF VALUE. OTHER PARAMETERS ARE
 C N = NUMBER OF DATA POINTS
 C IPAK = 0 INDICATES USER WISHES THE COMPUTER TO USE THE LARGEST A(X) PEAKS IN DESCENDING ORDER. (MAX OF 10 PEAKS)
 C = ANY NUMBER > 16: USER WISHES THIS NUMBER OF PEAKS TO BE USED. (IN DESCENDING ORDER - MAX OF 15 PEAKS)
 C AX(N) = REAL ARRAY CONTAINING THE A(X) VALUES
 C OTHER PARAMETERS USED IN THE SUBROUTINE
 C ZMAX = ESTIMATE OF MAXIMUM DEPTH TO MAGNETIC SOURCE
 C K = NUMBER OF PARAMETERS (3 * NUMBER OF SYMMETRIC COMPONENTS)
 C B(?) = REAL ARRAY CONTAINING THE PARAMETER VALUES
 C KMAX = MAXIMUM NUMBER OF PARAMETERS ALLOWED BY DIMENSION STATEMENTS
 C (MUST BE DIVISIBLE BY THREE)
 C A MAXIMUM OF 15 PEAKS (45 PARAMETERS) IS ALLOWED FOR THE LEAST SQUARE ROUTINE(MARQ1)
 C MAXIMUM COMMON USAGE: X(N),Q(2*KMAX),T(K/3),XX(0),XX(0),B(KMAX)
 COMMON X(200),Q(200),T(200),XX(600)
 COMMON /RESID/ZMAX,ICALL
 COMMON /SYM/B(45),K
 COMMON /CUT/CUTX,CUTRS,DAMIN,AMAX
 DIMENSION AX(1),IDIST(50),IAX(50)
 C ----FORMAT STATEMENTS----
 100 FORMAT(1H1,'PARAMETERS USED FOR THE DECOMPOSITION OF THE A(X)=',
 '\$! T(X)*2+T1(X)*#2 CURVE INTO ITS SYMMETRIC COMPONENTS; ',//,
 '\$! MAXIMUM VALUE OF A(X) = ',1PE14.6,' ! MINIMUM VALUE OF A(X) = ',
 '\$! 1PE14.6,' ! CUT OFF VALUE FOR A(X) = ',1PE14.6,' ! NUMBER OF
 '\$! SYMMETRIC PEAKS WITH A(X) VALUES LARGER THAN THE CUTOFF = !'
 '\$! 15,' ! NUMBER OF PEAKS REQUESTED = ',15,' ! TOTAL NUMBER OF !'
 '\$! PEAKS DETERMINED = ',15,' ! NUMBER OF PEAKS USED FOR LEAST !'
 '\$! SQUARE FIT = ',15,' /)
 101 FORMAT(1H1,'THE LOCATION AND A(X) VALUES FOR ALL PEAKS DETERMINED
 '\$ARE: ',/,' ! X - POSITION A(X) VALUE ',/(F9.3,10X,E15.7))
 102 FORMAT(1H1,'PARAMETERS OF THE SYMMETRIC CURVES IN A(X): (USED AS
 '\$ THE INITIAL VALUES IN THE LEAST SQUARE FIT PROCESS)',/,'10X,
 '\$ ALPHA',4X,'DEPTH(HOURS)',4X,'DIST(HOURS)',7X,'A(X) VALUE')

3 LEVEL 20

SYMPR3

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```
103 FORMAT(1H ,(3F15.3,F19.7))
104 FORMAT(//,* NO PEAKS WERE FOUND IN THE A(X) CURVE*)
105 FORMAT(//,* ***WARNING*** NO SYMMETRIC PEAKS HAVE A(X) VALUES
$> CUTOFF --- ANALYSIS ALLOWED TO CONTINUE*)
KMAX=45
L=2
M=N-1
LL=1
KL=0
DX=X(2)-X(1)
C      -----FIND LOCAL MAXIMA-----
DO 1 I=L,M
IF(AX(I).LE.AX(I+1).OR.AX(I).LE.AX(I-1)) GO TO 1
C      -----CHECK 2ND DERIVITIVE CRITERIA-----
DER2=ABS((AX(I+1)-2.*AX(I)+AX(I-1))
DEP1=SQRT((2.*AX(I)/DER2)*DX
IF(DEP1.GT.ZMAX) GO TO 1
IF(LL.GT.KMAX) GO TO 1
C      -----STORE PEAK LOCATIONS (INCREASING X-DIST-----
IF(AX(I).GE.CUTAX) KL=KL+1
IDIST(LL)=I
IF(LL.NE.1) GO TO 3
IAX(I)=I
GO TO 10
3 KK=1
C      -----STORE PEAK LOCATIONS (DECREASING VALUES OF AX(I)-----
4 IT=IAX(KK)
IF(AX(I).GT.AX(IT)) GO TO 5
KK=KK+1
IF((KK.GE.LL) GO TO 6
GO TO 4
6 IAX(KK)=I
GO TO 10
C      -----REARRANGE ARRAY IAX-----
5 J=KK
JJ=J
IM=IAX(J)
7 JJ=JJ+1
IF(JJ.EQ.LL) GO TO 8
IL=IAX(JJ)
IAX(JJ)=IM
IM=IL
GO TO 7
```

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8 IAX(JJ)=IM

IAX(J)=1

10 LL=LL+1

-----DETERMINE NO. OF PEAKS TO USE-----

1 CONTINUE

LL=LL-1

IF(IPEAK.NE.0) GO TO 9

KUSE=KL

IF(KL.GT.10) KUSE=10

GO TO 11

9 KUSE=IPEAK

IF(KUSE.GT.LL) KUSE=LL

IFIPEAK.GT.KMAX/3) KUSE=KMAX/3

11 CONTINUE

WRITE(6,100) AMAX,AMIN,CUTAX,KL,IPEAK,LL,KUSE

NO=-1

IF(LL.EQ.0) GO TO 50

-----FIND PARAMETERS (INCREASING X-DIST)-----

DO 20 J=1,LL

NO=NO+2

I=IDIST(J)

Q(NO)=X(I)

Q(NO+1)=AX(I)

20 CONTINUE

NO=NO+1

WRITE(6,101) Q(I),I=1,NO)

NO=-2

IF(KUSE.NE.0) GO TO 21

WRITE(6,105)

KUSE=1

21 CONTINUE

-----FIND PARAMETERS (DECREASING AX(I) VALUES-----

DO 30 J=1,KUSE

NO=NO+3

I=IAX(J)

C -----COMPUTE INITIAL DEPTHS-----

TEM=DEPIN(I,AX)

B(NO)=SQRT(TEM**2*AX(I))

B(NO+1)=TEM

B(NO+2)=X(I)

T(J)=AX(I)

30 CONTINUE

WRITE(6,102)

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```
L=-2
DO 40 J=1,KUSE
L=L+3
M=L+2
WRITE(6,103) B(I),I=L,M),T(J)
40 CONTINUE
C -----FINAL NU. OF PARAMETERS USED-----
K=KUSE*3
RETURN
50 WRITE(6,104)
K=0
RETURN
END
```

LEVEL 20

CORLAT DATE = 73332

17/28/02

SUBROUTINE CORLAT(N,A,B,COR)

C THIS SUBROUTINE COMPUTES THE PRODUCT-MOMENT CORRELATION COEFFICIENT BETWEEN
C TWO DATA ARRAYS A(N) AND B(N). THE MATHEMATICAL EXPRESSION IS:

$$COR = (\sum A^*B)/\sqrt{(\sum A^*A)(\sum B^*B)}$$

DIMENSION A(1),B(1)

REAL*8 X,Y,Z

X=0.0D0

Y=0.0D0

Z=0.0D0

DO 1 I=1,N

X=X+A(I)*B(I)

Y=Y+A(I)**2

1 Z=Z+B(I)**2

-----CORRELATION COEFFICIENT-----

COR=DABS(X/DSQRT(Y*Z))

RETURN

END

6 LEVEL 20

DATE = 73332

17/28/02

DEP

```
FUNCTION DEP(N,I,AX)
C THIS FUNCTION ESTIMATES THE INITIAL DEPTHS TO THE MAGNETIC SOURCE BY
C FINDING THE 1/2 WIDTH OF THE SYMMETRIC PEAKS IN THE AX(N) CURVE
C N = NO. OF DATA POINTS IN A(N)
C I = THE SUBSCRIPT OF THE ARRAY A(I) AT WHICH THE INITIAL DEPTH IS
C TO BE ESTIMATED
C AX(N)= A REAL ARRAY CONTAINING THE DATA CURVE
C MAXIMUM COMMON USAGE: X(N),XX(0)
COMMON X(200),XX(1000)
DIMENSION AX(1)
DX=X(2)-X(1)
JK=I-1
IF(N-I.LT.1) JK=N-1
LO=1
DO 21 JL=1,JK
IF(LO.EQ.JK) GO TO 25
IF(AX(I+1+LO).GT.AX(I+LO).OR.AX(I-1-LO).GT.AX(I-LO)) GO TO 25
LO=LO+1
21 CONTINUE
C -----COMPUTE 1/2 WIDTH-----
25 DEP=(DX*FLOAT(LO))/2.
RETURN
END
```

6 LEVEL 20

PEAKFT

DATE = 73332

17/28/02

SUBROUTINE PEAKFT(N,AX)

C THIS SUBROUTINE GIVES A QUANTITATIVE ESTIMATE AS TO THE QUALITY OF THE
C FIT BETWEEN THE A(X) CURVE AND THE COMPUTED CURVE - FOR EACH SYMMETRIC
C PEAK SEPARATELY. THE ANALYSIS FOR EACH PEAK IS DONE BETWEEN:

C PEAK POSITION +/- 3.*COMPUTED DEPTH
C AT THESE DISTANCES, THE CURVE FOR THE SYMMETRIC PEAK SHOULD (THEORETICALLY)
C BE AT 10% OF ITS MAXIMUM VALUE. THE PARAMETERS USED ARE:

C N=NQ. OF DATA POINTS IN THE AX(N) CURVE
C AX(N)=REAL ARRAY CONTAINING THE A(X) CURVE
C MAXIMUM COMMON USAGE: X(N),Y(0),C(0),D(0),E(0),DIFF(N),B(KMAX)
C F(0),AA(100),BB(100),XX(0)

REAL AA,BB

REAL MARQIP

COMMON X(100),Y(200),C(200),D(200),DIFF(200),XX(200)

COMMON /SYM/B(45),K

COMMON /PRMCRV/F(200)

COMMON /WORK/ AA(100),BB(100)

DIMENSION AX(1)

100 FORMAT(//1,ESTIMATION OF THE "GOODNESS OF FIT" FOR EACH SYMMETRIC

\$ PEAK - BASED UPON THE COMPARISON OF THE A(X) CURVE AND THE //,

\$ COMPUTED PARAMETER CURVE IN THE VICINITY OF THE SYMMETRIC PEAK!

\$ //, PEAK LOCATION(HRS) COMPARISON PTS FROM //,

\$ TO SUM OF SQUARES OF RESIDUALS CORRELATION COEFFICIENT

\$NT!)

102 FORMAT(6X,F6.3,T27,13,T51,F5.2,2X,F5.2,T73,E16.7,T103,E16.7)

WRITE(6,100)

DX=X(2)-X(1)

DO 1 I=1,K,3

TDEP=3.*B(I+1)

DISP=B(I+2)-X(1)

BDIST=(DISP-TDEP)/DX

FDIST=(DISP+TDEP)/DX

IA=IFIX(FDIST)+2

IB=IFIX(BDIST)+1

IFI(IA.GT.N) IA=N

IF(TB.LT.1) IB=1

NPT=IA-IB+1

IF(NPT.LT.100) GO TO 4

NR=(NPT-99)/2

IA=IA-NR

IB=IB-NR

NPT=IA-IB+1

4 DO 2 J=1,NP1

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PEAKFT

```
1 AA(J)=AX(IB+J-1)
2 BB(J)=F(IB+J-1)
CALL CORLAT(NPT,AA,BB,COR)
DO 3 J=1,NPT
3 AA(J)=DIFF(IR+J-1)
SQR=MARQIP(AA,AA,NPT)
WRITE(6,102) B(I+2),NPT,X(IR),X(IA),SQR,COR
1 CONTINUE
RETURN
END
```

LEVEL 20

DATE = 73332

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```

SUBROUTINE SETUP(N,TMAG,V,YMAX,YMIN)
C THIS SUBROUTINE SETS UP THE SCALE FACTORS AND MODIFIES TMAG(N) SO THAT THE
C TOTAL FIELD MAGNETICS CAN BE PLOTTED WITH THE /A(X)/*2 CURVE. THE SCALE
C IS SET SUCH THAT 25 SMALL DIVISIONS ON THE Y-AXIS =1000 GAMMA (2000 GAMMA
C IN SPECIAL CASES). V(N) CONTAINS THE THE MODIFIED TMAG VALUES. YMIN IS
C THE MINIMUM VALUE OF THE /A(X)/*2 CURVE, YMAX IS THE MAXIMUM VALUE.
DIMENSION TMAG(1),V(1)
CALL MINMAX(N,TMAG,TMIN,TMAX)
CS=(YMAX-YMIN)/4000.
CP=(YMAX+YMIN)/2.
IF((TMAX-TMIN).GT.2000.) GO TO 2
DO 1 I=1,N
 1 V(I)=((TMAG(I)-TMIN)*CS)+CP
      RETURN
2 CS=0.5*CS
DO 3 I=1,N
 3 V(I)=((TMAG(I)-TMIN)*CS)+CP
      RETURN
END

```

LEVEL 20

MINMAX

DATE = 73332

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SUBROUTINE MINMAX(N,X,XMIN,XMAX)

C THIS SUBROUTINE FINDS THE MINIMUM AND MAXIMUM VALUES WITHIN A 1-DIMENSIONAL

ARRAY X(N)

DIMENSION X(1)

XMIN=X(1)

XMAX=X(1)

DO 10 I=2,N

IF(X(I).GT.XMAX) XMAX=X(I)

IF(X(I).LT.XMIN) XMIN=X(I)

10 CONTINUE

RETURN

END

G LEVEL 20

PLOT DATE = 73332

17/28/02

```
SUBROUTINE PLOT(N,NCUR,U,V,W)
C THIS WILL PLOT UP TO A MAXIMUM OF THREE(3) CURVES: U(N),V(N),W(N) WITH A
C COMMON X-AXIS GIVEN BY X(N). THE ARRAY X(N) MUST BE STORED IN COMMON
C CURVES ARE PLOTTED USING SYMBOLS + * . RESPECTIVELY
C NCUR = NO. OF CURVES TO BE PLOTTED
C THIS SUBROUTINE REQUIRES THE USE OF SUBROUTINE PRLPLT
C MAXIMUM COMMON USAGE: X(N),XX(0)
COMMON X(200),XX(1000)
DIMENSION U(1),V(1),W(1),Y(200,3),SYM(3)
DATA SYM(1)'+'/SYM(2)'*'//SYM(3)'.'/
C -----SETUP PLOT ARRAYS-----
2 DO 1 I=1,N
  Y(I,1)=U(I)
  Y(I,2)=V(I)
  1 Y(I,3)=W(I)
C -----FIND SMALLEST AND LARGEST VALUE-----
10 CALL MINMAX(N,U,YMIN,YMAX)
CALL MINMAX(N,V,XMIN,XMAX)
CALL MINMAX(N,W,ZYMIN,ZYMAX)
CALL MINMAX(N,X,XMIN,XMAX)
YMIN=AMIN(YMIN,XMIN,ZYMIN)
YMAX=AMAX(YMAX,XMAX,ZYMAX)
C -----PLOT THE CURVES-----
CALL PRLPLT(XMAX,XMIN,YMAX,YMIN,NCUR,N,N,SYM,Y,X)
RETURN
END
```

6 LEVEL 20

DATE = 73332

17/28/02

PRLPLT

SUBROUTINE PRLPLT(XMAX,XMIN,YMAX,YMIN,LINES,LAST,SYM,X)

THIS SUBROUTINE IS A PLOT ROUTINE WHICH HAS THE PARAMETERS:

NY = NUMBER OF PLOTS TO BE MADE (FOR EACH PLOT A SYMBOL MUST
BE ASSIGNED IN THE CALLING PROGRAM - EX. DATA SYM(1)/* /

LINES = THE TOTAL NUMBER OF LINES ALONG THE X AXIS TO BE PLOTTED

LAST = THE NUMBER OF THE LAST POINT TO BE PLOTTED

SYM = AN ARRAY CONTAINING THE SYMBOLS TO BE USED IN PLOTTING

Y = A 2-DIMENSIONAL ARRAY CONTAINING THE Y VALUES TO BE PLOTTED

X = A 1-DIMENSIONAL ARRAY CONTAINING THE X AXIS VALUES

DATA BLANK// !,CR/-!,PL/+!/

DIMENSION SYM(NY),ZY(11),GRAPH(101),Y(200,3),X(1)

15 FORMAT(//,8X,P11E10.2)

20 FORMAT(1H,3X,'X VALUES',2X,20('+'1.5X,1Y VALUES'))

75 FORMAT(1H,2X,1PE9.2,2X,101A1,2(1PE9.2))

76 FORMAT(1H,2X,1PE9.2,2X,101A1,2X,1PE11.4)

85 FORMAT(1H,8X,P11E10.2)

90 FORMAT(1HO,3X,'SCALE = ',1PE11.4,4X,'XSCALE = ',1PE11.4)

XSCALE=(XMAX-XMIN)/(LINES-1.)

YSCALE=(YMAX-YMIN)/100.

DO 10 K=1,11

10 ZY(K)=10.*(K-1)*YSCALE+YMIN

WRITE(6,15) (ZY(K),K=1,11)

WRITE(6,20)

INDEX=1

40 X1=XMIN

DO 80 I=1,LINES

IF(MOD(I,5).EQ.1) GO TO 45

GRAPH(1)=CR

GRAPH(101)=CR

GO TO 50

45 GRAPH(1)=PL

GRAPH(101)=PL

50 DO 55 J=2,100

55 GRAPH(J)=BLANK

57 DO 60 J=1,NY

IY=(Y(INDEX,J)-YMIN)/YSCALE+1.5

GRAPH(IY)=SYM(J)

60 CONTINUE

INDEX=INDEX+1

60 TO 70

70 IF(NY.EQ.1) GO TO 71

WRITE(6,75)X(1),(GRAPH(J),J=1,101),Y(1,1),Y(1,2)

60 TO 80

PRLPL019

PRLPL020

PRLPL021

PRLPL022

PRLPL023

PRLPL024

PRLPL025

PRLPL026

PRLPL027

PRLPL028

PRLPL029

PRLPL030

PRLPL031

PRLPL032

PRLPL033

PRLPL034

PRLPL035

PRLPL036

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PRLPL039

PRLPL040

PRLPL041

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PRLPL045

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PRLPL053

PRLPL054

PRLPL055

PRLPL056

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PRLPLT

```
71 WRITE(6,76)X(I),(GRAPH(J),J=1,101),Y(I,1)
80 CONTINUE
      WRITE(6,20)
      WRITE(6,85) (ZY(K),K=1,11)
      WRITE(6,90) YSCALE,XSCALE
      RETURN
END
```

PRLPL063
PRLPL064
PRLPL065
PRLPL067
PRLPL069

LEVEL 20

MARQ1 DA

17/28/02

SUBROUTINE MARQ1(N,Y1*)

MODIFICATION OF STANFORD UNIVERSITY
LIBRARY PROGRAM NUMBER C008
WILLIAM E RIDDLER (SSCC)
MARCH 9, 1967 REVISED APRIL 4, 1968

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C MODIFICATION OF STANFORD UNIVERSITY
C LIBRARY PROGRAM NUMBER C008
C WILLIAM E RIDDLE (SCC)
C MARCH 9, 1967 REVISED APRIL 4, 1968

C THIS SUBROUTINE DOES A NONLINEAR LEAST SQUARE ANALYSIS BASED ON THE METHOD
C GIVEN BY MARQUARDT. THE ROUTINE FINDS THE BEST FIT FOR THE K PARAMETERS
C (OR K/3 SYMMETRIC CURVES) OF THE AX(N) CURVE. THE INITIAL VALUES OF THE
C PARAMETERS (FOUND BY SYMPR3) ARE IMPROVED BY MINIMIZING THE SUM OF THE
C SQUARES OF THE DIFFERENCE BETWEEN THE AX(N) CURVE AND THE LEAST SQUARE
C CURVE. THE EFFECTIVENESS OF THE METHOD IS HIGHLY DEPENDENT UPON THE
C QUALITY OF THE INITIAL VALUES FOR THE PARAMETERS.
C FOR EACH ITERATION, THE 1ST PARTIAL DERIVATIVES OF ALL PARAMETERS ARE
C COMPUTED ONCE
C FOR EACH FUNCTION EVALUATION, THE PARAMETER VALUES ARE USED TO COMPUTE THE
C LEAST SQUARE APPROXIMATION CURVE.
C MAXIMUM COMMON USAGE: X(N),TM(N),SS(5*KMAX/3),TEM(0),DIFF(N),XX(0)
C B(KMAX),F(N)

C SUBROUTINE PARAMETERS
INTEGER N,K,IMAX
REAL LAMRDP,NU,EPS,P,B,Y,F

C LOCAL VARIABLES
REAL TAU,FACT,SUM,S,NORMD,NORMG,DIFF,G,GS,GU,D,DS,DG,BI,A,AS
REAL MARQIP,LAMBDA
REAL LAMBDA1,LAMBDA2
INTEGER C1,C2,I,J,L,C3,II
DOUBLE PRECISION T1,T2,TSUM
COMMON X(200),TM(200),SS(200),TEM(200),DIFF(200),XX(200)
COMMON /LSTSQ/ LAMBDA,NU,EPS,OUTPT,SPD,IMAX,INPLT
COMMON /RESID/ ZMAX,ICALL
COMMON /SYN/ B(45),K
COMMON /PRMCRV/F(200)
COMMON /LSQERR/ IERR
DIMENSION P(200,45),Y(1)
LOGICAL OUTPT

DIMENSION G(45),GS(45),GU(45),DI(45),DS(45),DG(45),BI(45),
$ A(45,45),AS(45,45)

C SUPROGRAMS CALLED: MARQGS(SUBROUTINE), MARQIP(FUNCTION)

C -----FORMAT STATEMENTS----- MARQ1027
C -----FORMAT STATEMENTS----- MARQ1028

```

; LEVEL 20

MARQ1

DATE = 73332

17/28/02

```
$ SYMMETRIC COMPONENTS OF THE A(X) CURVE: ! /  
90 FORMAT(//,! GOODNESS OF FIT PRIOR TO LEAST SQUARE PROCESS: ! /,  
$! SUM OF SQUARES OF RESIDUALS=!,E16.7!,!  
$! CORRELATION COEFFICIENT=!,E18.7,/,! NUMBER OF ITERATIONS REQUESTED  
$TED=!,15)  
551 FORMAT(//,! THE LEAST SQUARE ROUTINE HAS CONVERGED TO A PROPER SOLUTION  
$ DURING ITERATION: ! 13, /)  
601 FORMAT(//! RESULTS FROM ITERATION=!,I4,/,! NO. OF FUNCTION EVALUATIONS  
$ONS=!,I4,/,! SUM OF SQUARES OF RESIDUALS=!,E16.8,/,! CORRELATION  
$ COEFFICIENT-A(X) AND COMPUTED CURVE=!,E17.7,/,! NO. OF CURVES REJECTED  
$ DURING THIS ITERATION: ! 14)  
901 FORMAT(//,! FINAL APPROXIMATION TO THE PARAMETERS OF THE SYMMETRIC  
$ CURVES - AFTER ! 13, ! ITERATIONS: ! ,/9X, ! ALPHA!,5X, ! DEPTH(HOURS)!,  
$4X, ! X-DIST(KM)!,4X, ! X-DIST(KM)!,/,!(3F15.5,X,  
$2F15.5))  
902 FORMAT(1H!,!THE VALUES FOR THE APPROXIMATION TO THE A(X) CURVE ARE  
$!: ! 10(1/,1P5E20.9))  
903 FORMAT(//,! -----ALL SYMMETRIC CURVES HAVE BEEN REJECTED DURING  
$ ITERATION: ! 13, ! OF THE LEAST SQUARE ALGORITHM-----! /,  
$! NO. OF FUNCTION EVALUATIONS=!,I4,/,! PARAMETERS REJECTED DURING  
$ THIS ITERATION: ! ,/!(3F15.5))  
904 FORMAT(//,! THE RESIDUAL VALUES (A(X) - COMPUTED) ARE: ! ,  
$10(/,1P5E20.9))  
905 FORMAT(1H!,! PLOT OF THE LEAST SQUARE APPROXIMATION TO THE A(X) C  
$URVE: (AFTER ! 13, ! CHECKS OF THE RESIDUAL VALUES) !  
$/,! (+) A(X) CURVE (*) LST SQ CURVE (!) RESIDUAL CURVE)  
906 FORMAT(! PARAMETERS OF REJECTED CURVES: ! ,!(3F15.5))  
907 FORMAT(! -----RECHECK THE A(X) CURVE AND THE INITIAL VALUES OF  
$ THE PARAMETERS-----! )  
908 FORMAT(//,! NOTE: THIS PLOT DOES NOT INCLUDE THE ! 13, ! SYMMETRIC  
$ CURVES WHICH HAVE BEEN REJECTED DURING THE LEAST SQUARE PROCESS.,!  
$/,! TO IMPROVE THE RESULTS REFER TO THE FOLLOWING EXPLANATION:  
$/!,! REJECTION CRITERIA! 27X, ! POSSIBLE CAUSE! 34X,  
$/!,! REMEDY! 1/, ! COMPUTED POSITION OUT OF DATA RANGE: ! ,T45, ! SMALL PEA  
$K NEAR OR ON LARGE PEAK OR ! T90, ! USE MORE PEAKS FOR ANALYSIS OR!  
$/!,! COMPUTED DEPTH NEGATIVE: ! ,T45, ! MANY LARGE PEAKS CLOSE TOGETHER  
$/!,! T90, ! USE ADDITIONAL RESIDUAL ANALYSIS: ! /, ! COMPUTED DEPTH LARGER  
$ THAN ZMAXKM: ! ,T45, ! ZMAXKM TOO SMALL OR LARGE ! ,T90, ! INCREASE ZMAXK  
$M!)  
909 FORMAT(//,! ***ERROR*** ! THE LEAST SQUARE SOLUTION HAS BLOWN  
$ UP. PROGRAM EXECUTION HAS TERMINATED ON AN EXPONENT OVERFLOW!  
$/,! OR UNDERFLOW. THE ERROR HAS BEEN CAUSED BY THE REJECTION OF AN  
$ IMPORTANT SYMMETRIC CURVE. THUS CAUSING THE LEAST SQUARE: ! /,
```

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MARQ1

C -----FILL ARRAY G(K)-----

TSUM=0. DO

DO 121 JJ=1,N

T1=DIFF(JJ)

T2=P(JJ,II)

121 TSUM=TSUM+T1*T2

120 G(II)=TSUM

C -----CONTINUE WITH ITERATIVE PROCEDURE-----

NORMG = MARQIP(G,K)

DO 150 I=1,K

150 DG(II) = 1.0 / SQRT(A(II,II))

DO 160 I=1,K

GS(II) = G(II) * DG(II)

DO 160 J=1,K

160 AS(II,J) = A(II,J) * DG(II) * DG(J)

FACT = 1.0

IF (LAMBDA*GT.(1.0E-12)) FACT = 1.0 / NU

C

200 DO 220 I=1,K

DO 210 J=1,K

210 A(I,J) = AS(I,J)

A(II,I) = AS(II,I) + LAMBDA * FACT

220 GU(II) = GS(II)

MN=K

CALL MARQGS (MM,A,GU,DS,\$1010)

DO 230 I=1,K

230 D(II) = DS(II) * DG(II)

C

300 DO 310 I=1,K

310 B1(II) = B(II) + D(II)

-----EVALUATE THE FUNCTION-----

CALL FUNCT(B1,K,F,N)

C1 = C1+1

DO 320 I=1,N

320 DIFF(II) = Y(II) - F(II)

S = MARQIP(DIFF,DIFF,N)

C

-----CHECK SUM OF SQUARES-----

IF (S-SUM) 400,400,500

C

-----SUM-----

MARQ1076

MARQ1077

MARQ1046

MARQ1047

MARQ1048

MARQ1049

MARQ1050

MARQ1051

MARQ1052

MARQ1053

MARQ1054

MARQ1055

MARQ1056

MARQ1057

MARQ1058

MARQ1059

MARQ1060

MARQ1061

MARQ1062

MARQ1063

MARQ1064

MARQ1065

MARQ1066

MARQ1067

MARQ1068

MARQ1069

MARQ1070

MARQ1071

MARQ1072

MARQ1073

MARQ1074

MARQ1075

MARQ1076

MARQ1077

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```

      '' SOLUTION TO BECOME UNSTABLE. REMEDY: CHANGE THE NECESSARY INPUT
      $ PARAMETERS (ZMAXM, SHPSPD) OR INPEAKT, NPEAKD, OR ICUTRSD) SO THAT THE CURVE IS NOT REJECTED.
      $/1. ZMAXM, SHPSPD - IF THE SYMMETRIC CURVE WAS REJECTED BECAUSE THE COMPUTED DEPTH(HRS) * SHPSPD) EXCEEDED ZMAXM. ! /, 5X,
      $/2. NPEAKT, NPEAKD - IF ERROR OCCURRED BEFORE A RESIDUAL ANALYSIS.
      $MORE SYMMETRIC CURVES ARE PROBABLY ! /, 26X, NEEDED FOR THE LEAST SQUARE SOLUTION. ! /, 5X, !3. CUTRSD - IF ERROR OCCURRED AFTER A RESIDUAL ANALYSIS. TOO MANY SYMMETRIC CURVES FOR CURVES TOO CLOSE TOGETHER ! /, 18X, HAVE BEEN ADDED. TRY LOWERING THE RESIDUAL PEAK CUTOFF LEVEL BY MAKING CUTRSD SMALLER. ! /, ! FOR HELP IN FINDING THE PROBLEM, THE FINAL VALUES OF THE PARAMETERS FOR ALL SYMMETRIC ! /, ! CURVES, AT THE TIME THE ERROR OCCURED ARE: (ALPHA, DEPTH(HRS),
      $X-POSITION(HRS)) ! , ! (3F15.5)

      IEX=0          MARQ1030
      IP=ICALL-1     MARQ1030
      TAU = 1.0E-3    MARQ1031
      C2 = 0          MARQ1033
      C1 = 1          MARQ1036
      SUM = MARQ1(I,DIFF,F,N)  MARQ1037
      C
      IF(INPLT.EQ.0) GO TO 70
      CALL PLOT(N,3,Y,F,TM)
      70 CALL CORLAT(N,Y,F,VAL)
      WRITE(6,90) SUM,VAL,TMAX
      WRITE(6,89)
      C
      C-----START ITERATIVE PROCEDURE-----
      C-----COMPUTE 1ST PARTIAL DERIVITIVES-----
      100 CALL DERIV(N,P)
      C
      NREJ=0          MARQ1040
      C2 = C2 + 1
      C3 = 0          MARQ1041
      DO 120 I=1,K    MARQ1042
      DO 110 J=1,K    MARQ1043
      C-----FILL ARRAY A(K,J)-----
      TSUM=0.D0
      DO 111 J=1,N
      T1=P(JJ,I)
      T2=P(JJ,J)
      111 TSUM=TSUM+T1*T2
      110 A(I,J)=TSUM

```

LEVEL 20 MARQ1 DATE = 73332 17/28/02 MARQ1078
 C -----EXIT FROM ROUTINE-----
 C 400 DO 410 I=1,K MARQ1079
 410 B(I) = B(I)
 LAMBDA = LAMBDA * FACT MARQ1080
 SUM = S
 GO TO 600 MARQ1081
 C
 C 500 IF (C3.NE.0) GO TO 550 MARQ1082
 C -----CHANGE CONVERGENCE FACTOR-----
 C NORMD = MARQ1P(D,K) MARQ1083
 IF (TERR.EQ.0) GO TO 125 MARQ1084
 WRITE(6,909) (B(I),I=1,K)
 STOP MARQ1085
 C
 125 CONTINUE MARQ1086
 S = NORMD * NORMG
 S = MARQ1P(D,G,K) / SQRT(S) MARQ1087
 IF (S.GT.(.70107)) GO TO 570 MARQ1088
 C -----CONTINUE WITH ROUTINE-----
 C
 FACT = FACT * NU MARQ1089
 GO TO 200 MARQ1090
 C
 C 550 IF (C3.NE.40) GO TO 570 MARQ1091
 WRITE(6,551) C2 MARQ1092
 GO TO 600 MARQ1093
 C
 C 570 C3 = C3 + 1 MARQ1094
 DO 580 I=1,K MARQ1095
 580 D(I) = D(I)/NU MARQ1096
 GO TO 300 MARQ1097
 C
 C -----CHECK PARAMETERS-----
 C 600 CALL PRMCHK(N,NREJ) MARQ1098
 IREJ=NREJ#3
 IF (NREJ.NE.0) IEX=IEX+NREJ
 IF (K.EQ.0) GO TO 700
 CALL CORLAT(N,Y,F,VAL)

3 LEVEL 20

MARQ1

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C -----PRINT RESULTS-----

WRITE(6,601) C2,C1,SUM,VAL,NREJ
IF(IREJ.NE.0) WRITE(6,906) (SS(I),I=1,IREJ)

602 IF(C2.EQ.1MAX) GO TO 900

IF(VAL.LT.0.98) GO TO 100

900 J=1

DO 910 I=1,K,3

DO 911 J=1,3

SS(J)=B(I+JJ-1)

911 J=J+1

DO 910 JJ=1,2

SS(J)=B(I+JJ)*SPD

910 J=J+1

KK=5*K/3

WRITE(6,901) C2,(SS(I),I=1,KK)

IF(NREJ.EQ.0) GO TO 915

CALL FUNCT(R,K,F,N)

915 CALL PEAKFT(N,Y)

IF(.NOT.OUTPUT) GO TO 710

WRITE(6,902) (F(I),I=1,N)

WRITE(6,904) (DIFF(I),I=1,N)

C MARQ1136

C MARQ1137

C 710 WRITE(6,905) IP

C -----PLOT RESULTS-----

CALL PLOT(N,3,Y,F,DIFF)

IF(IEX.EQ.0) GO TO 1000

WRITE(6,908) IEX

GO TO 1000

700 WRITE(6,903) C2,C1,(SS(I),I=1,IREJ)

WRITE(6,907)

1000 RETURN

1010 RETURN 1

C LAST CARD OF MARQ1

END

MARQ1138

MARQ1139

MARQ1140

LEVEL 20 MARQGS DATE = 73332

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```

SUBROUTINE MARQGS(M,A,B,X,*)
C CALLED BY MARQ1 - LIBRARY PROGRAM C 008      MARQG000
C SOLVES, BY GAUSSIAN ELIMINATION, THE SET OF SIMUL. EQUATIONS. A*X=B   MARQG001
C PARAMETERS                                         MARQG002
C
C      INTEGER M                                     MARQG003
C      REAL A,B,X                                     MARQG004
C      DIMENSION A(45,45),B(1),X(1)                 MARQG005
C
C LOCAL VARIABLES                               MARQG006
C      INTEGER I,J,IMAX,K,L,M                      MARQG008
C      REAL MX,T,QUOT                                MARQG009
C
C
C      DO 500 K= 1,M                                MARQG010
C      MX = 0.0                                      MARQG011
C      IMAX = K                                     MARQG012
C      DO 100 L = K,M                               MARQG013
C      IF (MX.GT.*ABS(A(I,K))) GO TO 100          MARQG014
C      MX = ABS(A(I,K))                           MARQG015
C      IMAX = I                                     MARQG016
C
C      100 CONTINUE                                 MARQG017
C      IF (MX.EQ.0.0) GO TO 1010                  MARQG018
C      I = K                                       MARQG019
C      IF (K.EQ. IMAX) GO TO 400                  MARQG020
C
C      MX = ABS(A(I,K))                           MARQG021
C
C      J=IMAX                                     MARQG022
C      T=B(I)                                     MARQG023
C      B(I)=B(J)                                  MARQG024
C      B(J)=T                                     MARQG025
C
C      DO 200 L=1,M                                MARQG026
C      T = A(I,L)                                 MARQG027
C      A(I,L) = A(J,L)                            MARQG028
C      200 A(J,L) = T                            MARQG029
C
C
C      400 L=K+1                                 MARQG030
C      IF(L.GT.*M) GO TO 500                      MARQG031
C      DO 450 J=L,M                               MARQG032
C      QUOT = A(J,K) / A(I,K)                   MARQG033
C      B(J) = B(J) - B(K)*QUOT                  MARQG034
C      DO 450 I=L,M                               MARQG035
C      A(J,I) = A(J,I) - QUOT*A(K,I)            MARQG036
C
C
C      450 A(J,I) = A(J,I) - QUOT*A(K,I)          MARQG037
C
C
C

```

LEVEL 20

MARQGS

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C MARQG042
C MARQG043
C MARQG044
C MARQG045
C MARQG046
C MARQG047
C MARQG048
C MARQG049
C 500 CONTINUE
C DO 590 K=1,M
I = M + 1 - K
T = 0
J = I+1
IF(J.GT.M) GO TO 590
DO 580 L=J,M
580 T = T + A(I,L) * X(L)
590 X(I) = (B(I)-T)/A(I,I)
C RETURN
1010 RETURN 1
END

LEVEL 20

MARQIP DATE = 73332

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```
FUNCTION MARQIP (V1,V2,N)
C CALLED BY MARQ1 - LIBRARY PROGRAM NUMBER C 008
C CALCULATES DOUBLE PRECISION INNER PRODUCT
C
C PARAMETERS
C COMMON /LSQERR/ IERR
C INTEGER N
C REAL V1, V2, MARQIP
C DIMENSION V1(N), V2(N)
C LOCAL VARIABLES
C INTEGER I
C DOUBLE PRECISION T1, T2, SUM
C
C SUM=0.D0
DO 10 I=1,N
T1 = V1(I)
T2 = V2(I)
IF(T1*T2.LT.1.E60.AND.T1*T2.GT.1.E-60) GO TO 8
IERR=1
RETURN
8 CONTINUE
10 SUM = SUM + T1 * T2
MARQIP = SUM
RETURN
END
```

LEVEL 20 DERIV

DATE = 73332 17/28/02

SUBROUTINE DERIV(N,P)

C THIS SUBROUTINE IS USED BY SUBROUTINE MARQ1 TO COMPUTE THE DERIVATIVES
C OF THE ANALYTIC EXPRESSION FOR THE SYMMETRIC BELL SHAPED CURVES WHICH
C COMPOSE THE AX(X) CURVE. THIS EXPRESSION IS F(X)=ALPHA**2/H**2+(X-PQS)**2
C N=NO. OF DATA POINTS
C P(N,K)= ARRAY CONTAINING THE FIRST PARTIAL DERIVATIVES OF THE FUNCTION
C OTHER PARAMETERS USED BY THE SUBROUTINE ARE:

C K= NO. OF PARAMETERS TO BE MINIMIZED

C B(K)= ARRAY CONTAINING THE PARAMETERS

C F(X) WITH RESPECT TO EACH OF THE PARAMETERS AND AT EACH DATA POINT

C MAXIMUM COMMON USAGE: X(N),XX(0),B(KMAX)

COMMON X(200),XX(1000)

COMMON /SYM/ B(45),K

REAL*8 D,C,CON,CON2

DIMENSION P(200,45)

L=K/3

DO 10 I=1,N

DO 10 J=1,K,3

DIST=X(I)

ALPHA=B(J)

DEPTH=B(J+1)

PKPOS=B(J+2)

C-----COMPUTE 1ST PARTIAL DERIVATIVES-----

D=2*D0*ALPHA**2

C=DIST-PKPOS

CON=DEPTH**2+C**2

CON2=CON**2

C-----ALPHA(A)-----

P(I,J)=D/(CON*ALPHA)

C-----DEPTH(H)-----

P(I,J+1)=-D*DEPTH/CON2

C-----POSITION(C)-----

10 P(I,J+2)=D*C/CON2

RETURN

END

LEVEL 20

FUNCTION

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```
C SUBROUTINE FUNCT(B,K,F,N)
C THIS SUBROUTINE IS USED BY SUBROUTINE MARQ1 TO DETERMINE THE VALUES OF THE
C FUNCTION (WHICH IS BEING USED FOR THE LEAST SQUARE MINIMIZATION) AT EACH
C OF THE DATA POINTS
C K=NO. OF PARAMETERS BEING MINIMIZED
C N= NO. OF DATA POINTS
C B(K)= ARRAY CONTAINING THE PARAMETER VALUES
C F(N)= ARRAY CONTAINING THE OUTPUT FUNCTION VALUES
C MAXIMUM COMMON USAGE: X(N),XX(0)
COMMON X(200),XX(1000)
REAL*8 SUM
DIMENSION B(1),F(1)
SUM=0.0D0
DO 3 I=1,N
  DO 2 J=1,K,3
    DIST=X(I)
    ALPHA=B(J)
    DEPTH=B(J+1)
    PKPOS=B(J+2)
C -----SUM OF ALL CURVES AT POINT X-----
 2  SUM=SUM+(ALPHA**2)/((DEPTH**2)+(DIST-PKPOS)**2)
    F(I)=SUM
 3  SUM=0.0D0
    RETURN
END
```

LEVEL 20

PRMCHK DATE = 73332

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SUBROUTINE PRMCHK(N,NREJ)

C THIS SUBROUTINE CHECKS THE PARAMETERS OF THE SYMMETRIC CURVES TO SEE IF:

- C 1. SOURCE DEPTHS ARE NEGATIVE
- C 2. SOURCE DEPTHS ARE GREATER THAN ZMAX
- C 3. X-POSITIONS LIE OUTSIDE THE RANGE OF THE DATA
C IF ANY CONDITION IS FOUND, THE ROUTINE THROWS AWAY THAT SYMMETRIC CURVE
C AND REARRANGES THE ARRAY CONTAINING THE PARAMETERS

C N=NO. OF DATA POINTS

C NREJ=NO. OF SYMMETRIC CURVES REJECTED

C OTHER PARAMETERS USED IN THE SUBROUTINE ARE:

C K=NO. OF PARAMETERS

C B(K)=REAL ARRAY CONTAINING THE PARAMETERS

C MAXIMUM COMMON USAGE: X(IN) T(I) T(MAX) XX(0)

C COMMON X(200),T(200),TM(200),XX(600)

C COMMON /RESID/ ZMAX,LICALL

C COMMON /SYM/ B(45),K

NT=1

L=K

C -----CHECK PARAMETERS-----

C DO 11 I=1,K,3

C IF(B(I+2).GT.X(N)).OR.B(I+2).LT.X(1)) GO TO 12

C IF(B(I+1).LT.0..OR.B(I+1).GT.ZMAX) GO TO 12

C GO TO 11

C 12 L=K-3*NT

C NP=3*NREJ

C NREJ=NREJ+1

C NT=NT+1

C -----REJECT BAD CURVES-----

C DO 13 J=1,3

C TM(NP+J)=B(I+J-1)

C 13 B(I+J-1)=0.

C 11 CONTINUE

C NT=1

C -----REARRANGE PARAMETER ARRAY-----

C DO 14 I=1,K

C IF(B(I).EQ.0.) GO TO 14

C B(NT)=B(I)

C NT=NT+1

C 14 CONTINUE

C K=L

C RETURN

C END

; LEVEL 20

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RSDCHK

SUBROUTINE RSDCHK(N, AX)

C THIS SUBROUTINE CHECKS THE RESIDUAL CURVE (IE AX(I)-COMPUTED) FOR
 C ADDITIONAL SYMMETRIC PEAKS WHICH DO NOT OCCUR AS LOCAL MAXIMA ON THE
 C A(X) CURVE. NEW PEAKS FOUND IN THE RESIDUAL CURVE (BY THE SAME CRITERIA
 C AS THE AX(I)) ARE ADDED TO THE LIST OF OLD SYMMETRIC PEAKS IF THEY ARE
 C LOCATED A DISTANCE $> 2 * \Delta X$ AWAY FROM AN OLD PEAK. (STATEMENTS 51 & 52)
 C NEW PEAKS ARE NOT KEPT IF THEY ARE LESS THAN THIS DISTANCE
 C N = NO. OF DATA POINTS IN A(X)

AX(N) = REAL ARRAY CONTAINING THE A(X) CURVE
 OTHER PARAMETERS USED IN THE SUBROUTINE ARE:

R(N) = REAL ARRAY CONTAINING THE RESIDUAL VALUES (A(X) - COMPUTED)

K = NO. OF PARAMETERS
 R(K) = REAL ARRAY CONTAINING THE PARAMETER VALUES

ZMAX = MAXIMUM ESTIMATED DEPTH TO THE MAGNETIC SOURCE BODY
 NADD = NO. OF CURVES ADDED FROM RESIDUAL ANALYSIS

ND = NO. OF CURVES DELETED BY THE RESIDUAL ANALYSIS
 KMAX = MAXIMUM NO. OF PARAMETERS ALLOWED

MAXIMUM COMMON USAGE: X(N), Q(0), REJ(KMAX), W(2*KMAX), R(N), F(KMAX)
 B(KMAX), FF(N)

COMMON X(200), Q(200), REJ(200), W(200), R(200), F(200)

COMMON /LSTSQ/ LAMBDA, NU, EPS, OUTPT, SPD, IMAX, INPLT

COMMON /RESID/ ZMAX, ICALL

COMMON /SYM/ B(45), K

COMMON /PRMCRV/ FF(200)

COMMON /CUT/ CUTAX, CUTRSD, AMIN, AMAX

DIMENSION AX(1)

FORMAT STATEMENTS--

100 FORMAT(/, !, -----, THE TOTAL NUMBER OF SYMMETRIC PEAKS HAS REACHED
 \$HED!, I3, !, -----, !)

101 FORMAT(IH1, !, RESULTS FROM ANALYSIS NUMBER!, I3, !, OF THE RESIDUAL CUR
 \$VE: (AFTER THE LEAST SQUARE PROCESS)!, I1, !, THE CUTOFF LEVEL FOR T

\$HE RESIDUAL PEAKS=!, 1PE14.6)

102 FORMAT(/, !, THE!, I3, !, NEW SYMMETRIC CURVES ADDED HAVE THE PARAMETER
 \$S: (ONLY RESIDUAL PEAKS WITH VALUES > CUTOFF ARE CONSIDERED)!, /,
 \$(3F15.5))

103 FORMAT(/, !, THE!, I3, !, OLD SYMMETRIC CURVES REMOVED HAVE THE PARAMET
 \$ERS!: /, (3F15.5))

104 FORMAT(/, !, THE!, I3, !, NEW PARAMETERS OF THE!, I3, !, SYMMETRIC CURVES
 \$(AFTER RESIDUAL ANALYSIS) ARE!: /, (3F15.5))

105 FORMAT(/, !, NO NEW SYMMETRIC CURVES HAVE BEEN ADDED!)

106 FORMAT(/, !, NO OLD SYMMETRIC CURVES HAVE BEEN REMOVED!)

107 FORMAT(/, !, THE VALUES OF THE PARAMETERS REMAIN UNCHANGED AFTER THE
 \$ RESIDUAL ANALYSIS!)

```

LEVEL 20          RSDCHK          DATE = 73332           17/28/02
108 FORMAT(1H1,'PLOT OF THE A(X) CURVE(+), NEW PARAMETER CURVE(*), AND
$ TOTMAG CURVE(.); (AFTER RESIDUAL ANALYSIS, I2, I, & BEFORE LST SQR
$ ANALYSIS',I2,I)
KMAX=45
NADD=0
NN=1
M=N-1
KN=K+1
DX=X(2)-X(1)
ND=0
WRITE(6,101) ICALL,CUTRSD
DO 10 I=1,K
10 W(I)=B(I)
DO 1 I=2,M
C      ---- FIND LOCAL MAXIMA -----
C      IF(R(I).LT.CUTRSD) GO TO 1
C      IF(R(I).LE.R(I+1).OR.R(I).LE.R(I-1)) GO TO 1
C      ---- CHECK 2ND DERIVITIVE CRITERIA -----
C      DER2=ABS(R(I+1)-2.*R(I)+R(I-1))
DEP1=SQRT(2.*R(I))/DER2)*DX
IF(DEP1.GT.ZMAX) GO TO 1
IF(KN-3*NOD.GT.KMAX) GO TO 7
NADD=NADD+1
C      ---- FIND INITIAL DEPTH -----
C      T=DEP(N,I,R)
C      ---- ADD TEMPORARY NEW PEAKS -----
F(NN)=SQRT(T**2*AX(I))
F(NN+1)=T
F(NN+2)=X(I)
C      ---- CHECK FOR REPETITIVE PEAKS -----
C      IF(NN.EQ.I) GO TO 4
C      ---- DELETE NEW PEAK IF WITHIN 2*DX OF OLD PEAK -----
2  DO 3 JJ=1,K13
51  IF(ABS(F(NN-1)-W(JJ+2)),GE,2.*DX) GO TO 3
     NADD=NADD-1
     GO TO 4
3  CONTINUE
C      ---- ADD NEW RESIDUAL PEAKS -----
DO 15 J=1,3
15 W(KN+J-1)=F(NN+J-4)
KN=KN+3
4  NN=NN+3
LL=1

```

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1 CONTINUE

IF(NADD.EQ.0) GO TO 8

C -----CHECK LAST TEMPORARY PEAK-----

DO 16 JJ=1,K,3

52 IF(ABS(F(NN-1)-W(JJ+2)).GE.2.*UX) GO TO 16

NADD=NADD-1

GO TO 8

16 CONTINUE

DO 17 J=1,I³

17 WIKN+J-1)=F(NN+J-4)

KN=KN+3

GO TO 8

7 KMAX=KMAX/3

WRITE(6,100) KMAX

8 NT=1

KR=ND*3

KL=K+1

C -----OUTPUT RESULTS-----

IF(NADD.NE.0) GO TO 20

WRITE(6,105)

GO TO 21

20 K=KN-1

WRITE(6,102) NADD,(W(I),I=KL,K)

21 IF(ND.NE.0) GO TO 22

WRITE(6,106)

IF(NADD+ND.NE.0) GO TO 23

WRITE(6,107)

GO TO 30

22 WRITE(6,103) ND,(REJ(I),I=1,KR)

23 CONTINUE

C -----REARRANGE THE PARAMETER ARRAY-----

DO 6 I=1,K

IF(W(I).EQ.0.) GO TO 6

B(NT)=W(I)

NT=NT+1

6 CONTINUE

K=K-3*ND

KL=K/3

C -----OUTPUT PARAMETER VALUES FOR NEW PEAKS-----

WRITE(6,104) K,KL,(B(I),I=1,K)

30 CONTINUE

IF(NADD+ND.EQ.0) GO TO 32

CALL FUNCT(B,KIFF,N)

RSDCHK

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```
DO 35 JT=1,N
 35 R(JT)=AX(JT)-FF(JT)
     CALL PEAKFT(N,AX)
 32 IF(INPLT.EQ.0) GO TO 31
     ITEM=2*ICALL
     WRITE(6,108) ICALL,ITEM
 31 ICALL=ICALL+1
     RETURN
END
```

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INPUT

SUBROUTINE INPUT(N,TIME,TMAG)

C THIS SUBROUTINE READS TIMES AND TOTAL FIELD MAGNETIC VALUES FROM CARDS

C (BARTLETT FORMAT) AND CONVERTS TIME INTO HOURS

DIMENSION TIME(11),TMAG(1)

100 FORMAT (3I6,F8.1,F2.0,F10.0)

200 FORMAT (/, ' DATA SUPPLIED BY THE USER: ')

201 FORMAT (1I6X,'YEAR DAY TIME',/, ' START PROFILE:', I4,I6,I5,

\$F4.1,/, ' END PROFILE:', 2I6,I5,F4.1)

TIME(1)=0.

N=1

READ(5,100) IY1, ID1, IH1, AM1, AS1, TMAG(1)

AM1=AM1+AS1/60.

1 READ(5,100,END=10) IY, ID, IH, AM, AS, TOTMAG

AM=AM+AS/60.

TEMP =TMDF(IY1, ID1, IH1, AM1, IY, ID, IH, AM)/60.

IF(TEMP.EQ.TIME(N)) GO TO 1

N=N+1

1 IF(N.GT.200) GO TO 10

TMAG(N)=TOTMAG

TIME(N)=TEMP

GO TO 1

10 WRITE(6,200)

WRITE(6,201) IY1, ID1, IH1, AM1, IY, ID, IH, AM

RETURN

ENTRY PRINT(SHR,EHR)

TEM=SHR+IH1

DO 11 I=1,2

IHR=TEM

AMI=(TEM-IHRI)*60.+AM1

ID1=ID1

IF(AMI.LT.60.) GO TO 12

AMI=AMI-60.

IHRI=IHR+1

12 IF(IHRI.LT.24) GO TO 13

IHRI=IHR-24

ID1=ID1+1

13 IF(I1.EQ.2) GO TO 11

ID1=ID1

IHS=IHR

AM1=AM1

TEM=EHR+IH1

11 CONTINUE

WRITE(6,201) IY1, ID1, IH1, AM1, IY, ID1, IHRI, AM1

Return

End

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TMDF

```
FUNCTION TMDF(IY1, ID1, IH1, AM1, IY2, ID2, IH2, AM2)
C THIS FUNCTION FINDS THE TIME DIFFERENCE IN MINUTES
TMDF = 1440*(ID2-ID1)+60*(IH2-IH1)+AM2-AM1
J=IY2-IY1
IF(J.EQ.0) RETURN
TMDF=TMDF+525600*j+1440*(IY2-1)/4-(IY1-1)/4
RETURN
END
```